

Calibration, event reconstruction, data analysis, and limit calculation for the LUX dark matter experiment

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The LUX experiment has performed searches for dark-matter particles scattering elastically on xenon nuclei, leading to stringent upper limits on the nuclear scattering cross sections for dark matter. Here, for results derived from 1.4×10^4 kg days of target exposure in 2013, details of the calibration, event-reconstruction, modeling, and statistical tests that underlie the results are presented. Detector performance is characterized, including measured efficiencies, stability of response, position resolution, and discrimination between electron- and nuclear-recoil populations. Models are developed for the drift field, optical properties, background populations, the electron- and nuclear-recoil responses, and the absolute rate of low-energy background events. Innovations in the analysis include *in situ* measurement of the photomultipliers' response to xenon scintillation photons, verification of fiducial mass with a low-energy internal calibration source, and new empirical models for low-energy signal yield based on large-sample, *in situ* calibrations.

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I. INTRODUCTION

The Lambda cold dark-matter model of the Universe is in excellent agreement with numerous different types of astrophysical observations [1,2]. According to the most recent results from Planck, the mass/energy content of the Universe comprises 26.8% cold dark matter (CDM), which is nearly a factor of 5.5 more than the 4.9% contribution of baryonic matter [3]. The identity of the CDM is unknown, but a number of theoretical models predict that dark-matter particles interact very weakly with ordinary matter. Such models include weakly interacting massive particles (WIMPs) [4,5] and asymmetric dark matter [6,7] and in both cases predict that the dark-matter particles can elastically scatter with nuclei, producing nuclear recoils with energies of order 1–100 keV. Such elastic scattering events could be detectable using sufficiently sensitive instruments. A positive direct detection of dark-matter interactions would open a window into physics beyond the Standard Model, give new insights to cosmology, and create a new field of direct dark-matter observations.

One of the most effective technologies for the direct detection of dark matter is the two-phase xenon (Xe) time projection chamber (TPC) [8–12]. Xenon has low intrinsic radiological backgrounds since it has no long-lived isotopes other than the extremely long-lived ^{136}Xe isotope, which undergoes double-beta decay with a half-life of 2.1×10^{21} yr [13,14] and, at the low energies relevant to WIMP detection, produces a background rate of events that is small in comparison to the electron scattering of solar

neutrinos [15]. Liquid xenon (LXe) is dense (2.9 g/cm^3) and produces large-scintillation light and charge signals from both electron-recoil (ER) and nuclear-recoil (NR) events. The Xe TPC has the ability to discriminate ER from NR on an event-by-event basis using the charge-to-light ratio and can be expanded to large homogeneous volumes. With excellent position resolution and short LXe gamma-ray and neutron scattering lengths, this technology allows excellent self-shielding and extremely low backgrounds at low energies.

The Large Underground Xenon (LUX) experiment [11] is a two-phase xenon-based dark-matter detector located at the 4850 ft level of the Sanford Underground Research Facility (SURF) [16] in Lead, South Dakota, USA. LUX was assembled and first operated in a dedicated surface facility at SURF starting in 2009. A follow-up test of LUX on the surface (Run 2) started in October 2011 and ended in February 2012 [17]. The Davis Campus, at a depth of 4850 ft, was built to house LUX and other experiments and was completed at the end of May 2012. LUX began installation in the Davis Campus in June 2012, and the cooldown of the detector began in January 2013. The first WIMP search and calibration data (Run 3) were acquired during the period March–October 2013. First results from Run 3 were announced in October 2013 and subsequently published in Physical Review Letters [18]. The data were shown to be consistent with a background-only hypothesis, allowing 90% confidence limits to be set on spin-independent WIMP-nucleon elastic scattering with a minimum upper limit on the cross section of $7.6 \times 10^{-46} \text{ cm}^2$ at a WIMP mass of $33 \text{ GeV}/c^2$, the most stringent constraint on the dark-matter scattering cross section at that time. Further extensive calibration studies, together with a large number of other improvements, enabled a more

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sensitive analysis of the Run 3 data to be performed, further improving the sensitivity, especially to lower mass WIMPs, and reducing the 90% minimum upper limit exclusion to $6 \times 10^{-46} \text{ cm}^2$, again for a WIMP mass of $33 \text{ GeV}/c^2$. This improved sensitivity was published as a second Physical Review Letter in early 2016 [19], with a further Letter presenting results for spin-dependent interactions [20]. Here, detailed and complete descriptions of the new calibrations are presented, together with the analysis methods that underpin these results, which were unable to be published in the short Letter format.

Recently, results from an extended 332-day exposure (Run 4) have been published [21], providing a further roughly four-fold improvement in sensitivity for high WIMP masses relative to previous results. With no evidence of WIMP nuclear recoils, WIMP-nucleon spin-independent cross sections above $2.2 \times 10^{-46} \text{ cm}^2$ are now excluded at the 90% confidence level. When combined with the previously reported LUX exposure, this exclusion strengthens to $1.1 \times 10^{-46} \text{ cm}^2$ at $50 \text{ GeV}/c^2$. Many of the innovations outlined here persist in the analysis of these most recent data.

II. LUX EXPERIMENT

LUX is a two-phase Xe TPC designed to detect both the prompt scintillation light and the delayed ionization electrons that result from ionizing radiation. A schematic of the LUX detector is shown in Fig. 1, and photographs of LUX without its inner vessel, within the time projection chamber, and mounted in its water tank shielding are shown in Figs. 2, 3, and 4.

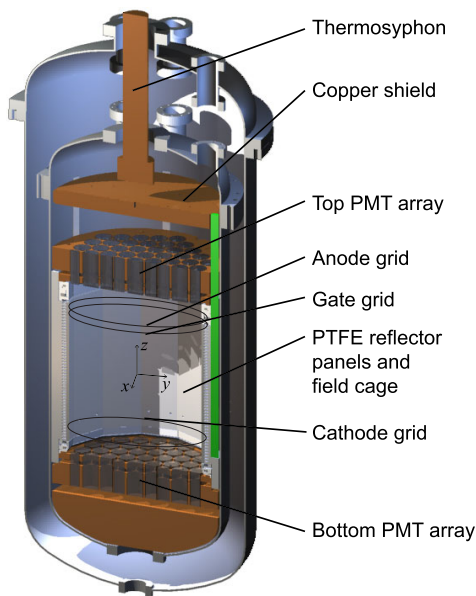


FIG. 1. Schematic of the LUX detector. LUX is a two-phase xenon time projection chamber containing an active LXe mass of 250 kg that is viewed from above and below by PMT arrays.

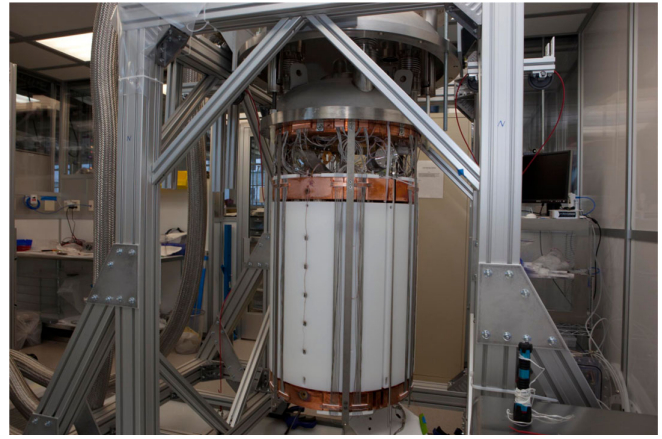


FIG. 2. Photograph of the LUX detector, with its inner vessel removed.

The central, fully active Xe region is defined by 12 tetrafluoroethylene (PTFE) reflector panels, the gate grid located 0.5 cm below the Xe liquid surface, and the cathode grid located 48 cm below the gate grid. The diameter is 47 cm. Field shaping rings, spaced by 2 cm, are mounted on the outer PTFE vessel and apply an electric field to drift the ionization electrons. An anode grid 0.5 cm above the surface, together with the gate grid, is used to generate high fields that extract the electrons and accelerate them, thus generating proportional scintillation light. Scintillation photons are produced almost uniformly throughout the gas gap, regardless of the electron extraction location. In Run 3, the LUX detector was operated with cathode, gate, and anode voltages of -10.0 , -1.5 , and $+3.5 \text{ kV}$, respectively. Finite-element methods were used to calculate field models that were found consistent with data [22,23]. The drift field at the detector center was 180 V/cm , with some position dependence (160 to 200 V/cm over the fiducial volume). The electron extraction field was $5.53 \pm 0.30 \text{ kV/cm}$ applied in the gas ($2.84 \pm 0.16 \text{ kV/cm}$ in the liquid), with the liquid level being the dominant uncertainty.

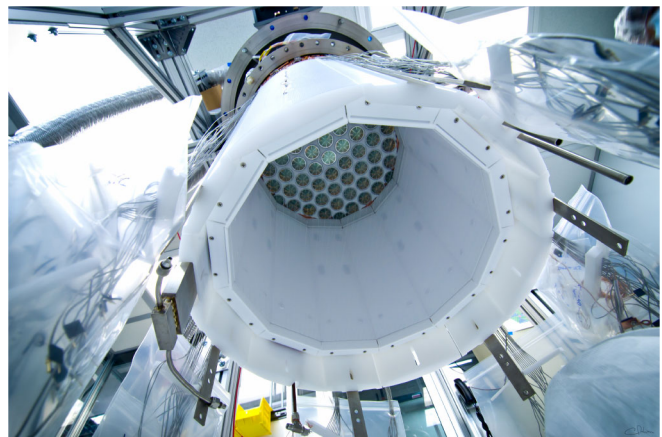


FIG. 3. Photograph of the inside of the LUX detector, with its lower PMT array removed.

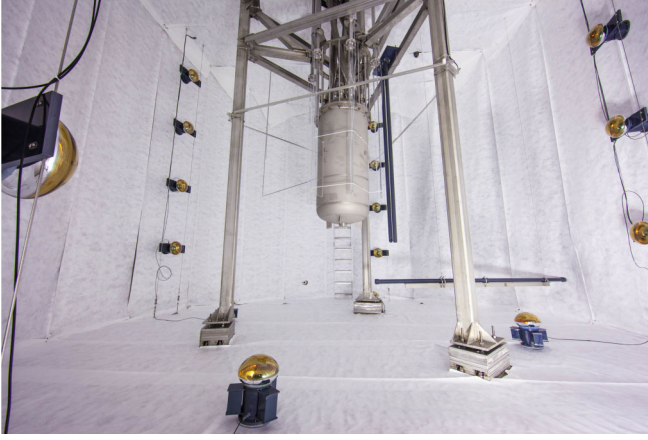


FIG. 4. Photograph of LUX installed in the Davis Cavern water shield at SURF.

An event in the LUX time projection chamber is characterized by two signals, corresponding to the detection of direct scintillation light (S1) and proportional light from ionization electrons (S2). The two light pulses occur within a maximum drift time of $322\ \mu\text{s}$, corresponding to a saturated electron drift velocity of $1.52\ \text{mm}/\mu\text{s}$ in LXe. Since electron diffusion in LXe is small, the S2 proportional scintillation pulse is produced in a small area at the gas-liquid interface. After corrections for any field nonuniformities, the original event may be located accurately in the xy -plane, allowing two-dimensional (2D) reconstruction. With precise z information from the drift time measurement, the three-dimensional (3D) event localization provides background discrimination via fiducial volume cuts.

Two arrays of 5.6 cm diameter Hamamatsu R8778 photomultiplier tubes (PMTs), 61 in each array, detect the S1 and S2 signals [24]. One PMT array above the liquid surface is primarily used to image the xy position of the proportional light pulse. The second PMT array is in the liquid, below the cathode grid. Liquid xenon has a high refractive index, 1.69 at 170 K, causing internal reflection at the liquid surface that in turn leads to most prompt light being collected in this bottom array. The high quantum efficiency of the R8778 PMTs, highly transparent grids, and use of PTFE reflectors between PMTs means that a very high light yield is achieved, measured to be 8.8 photoelectrons/keV electron-equivalent energy (hereafter keV_{ee}) for 122 keV γ rays at zero electric field.

The LUX Collaboration has introduced a number of innovations, including a low-radioactivity titanium cryostat, nitrogen thermosyphons, high-flow Xe purification, two-phase Xe heat exchangers, internal calibration with gaseous sources of $^{83\text{m}}\text{Kr}$ and ^3H , and nuclear-recoil calibration using multiple scatters of monoenergetic neutrons produced with a deuterium-deuterium (D-D) neutron generator. The LUX cryostat vessels were fabricated from Ti with very low levels of radioactivity [25], rivaling the purities achieved in copper. The two-phase TPC technique requires precise control of the

thermodynamic environment, and this was achieved in LUX through the development of a dedicated nitrogen thermosyphon system. This features precise, tunable, automated control delivering up to hundreds of kW of cooling power, plus reliable remote operation. A particularly important aspect of this system is that it has allowed highly controlled initial cooling of the LUX detector [11], which is necessary to avoid warping of the large plastic structures of the TPC. A related development is the purification system that allowed rapid circulation of Xe through an external gas-phase getter. Flow rates exceeding 27 standard L/min (229 kg/day) were achieved, while a stable liquid surface was maintained through the use of a weir. Negligible overall heat load on the detector was then obtained through the use of a two-phase heat exchanger system [26] and very efficient heat transfer between evaporating and condensing Xe streams. Xe purification was further monitored through the use of an innovative gas trapping and mass spectrometry system [27,28], sensitive to impurities at the sub-ppb concentration levels needed for good electron transport and light collection. This diagnostic capability allowed various portions of the gas system and detector to be monitored for contamination. Removal of Kr from the Xe, required to limit ^{85}Kr and ^{81}Kr beta-decay backgrounds, was performed before commencement of Run 3 using a charcoal column [29]. These systems were demonstrated during the experiment's first science run, in which cooldown was achieved in only 9 days. Sufficient LXe purity to begin science operations was achieved only one month after the initial filling with LXe, at which time the electron drift lifetime (the mean time an electron remains in the LXe before being absorbed by an impurity) was over $500\ \mu\text{s}$. Stable operation of the detector was maintained with mostly unattended operation over the five-month period, during which the pressure and liquid level had sufficient stability (1% and $< 500\ \mu\text{m}$, respectively) to introduce no measurable correlations in the S2 or S1 signals.

III. DATA ACQUISITION AND REDUCTION

A. DAQ configuration and single-photoelectron digitization acceptance

The signal of each PMT output is amplified by a factor of 5 with a linear preamplifier located at the instrumentation breakout and is subsequently shaped by a postamplifier that increases the pulse area by a further factor of 1.5. The postamplifier boards have additional amplification outputs to feed the LUX trigger and discriminator boards. The output of the postamplifier is digitized by 14-bit Struck SIS3301 analog-to-digital converter (ADC) boards at a sampling rate of 100 MHz (10 ns data samples). The Struck board firmware was modified to use pulse-only digitization (POD), a zero-suppression mode that only digitizes signals above a specified threshold. Since signals in LUX are dominated by long periods of baseline with short bursts of S1 (width $< 100\ \text{ns}$) and S2 (width $< 5\ \mu\text{s}$) signals, the

POD mode significantly reduces the number of recorded samples while increasing the maximum allowed acquisition rate. The signal threshold to begin digitizing the PMT signal was set to 1.5 mV at the Struck input; the signal threshold to end digitizing was set to 0.5 mV. An additional 24 samples before and 31 samples after the threshold crossings were also digitized. The LUX analog signal chain and data acquisition system (DAQ) maintain linearity in S1 and S2 signals at energies of ~ 100 keV_{ee}, well above the WIMP region of interest and comfortably above the ^{83}mKr and tritium ER calibrations. For a more detailed description of the LUX DAQ system, refer to Ref. [30].

PMT signals are continuously recorded by the DAQ regardless of trigger conditions, and the trigger pulse is digitized as an additional DAQ channel. Offline software, called the Event Builder, subsequently matches PMT signals within a specified time window around the trigger pulse for data processing and analysis. In this way, trigger changes can be made offline with no data loss.

For the nominal PMT gains of 4×10^6 , a single photoelectron (sphe) generates a pulse with 4 mV amplitude and full width at half maximum (FWHM) of ~ 20 ns. A typical sphe pulse is shown in Fig. 5 with height markers for the mean sphe height (μ_{sphe}) and its standard deviation ($\mu_{\text{sphe}} \pm \sigma_{\text{sphe}}$). The total noise from the electronics chain and the ADC, as measured at the Struck input, is $155 \mu\text{V}_{\text{rms}}$. The 1.5 mV POD threshold is indicated in Fig. 5, as is the height of a 5σ noise fluctuation.

To estimate the sphe digitization acceptance for each PMT, η_{sphe} , a Gaussian distribution, truncated at the POD threshold of 1.5 mV, was fitted to each sphe spectrum. This sphe calibration was performed with 400 nm light pulses from the light-emitting diode (LED) calibration system, which does not cause double-photoelectron emission and therefore results in values for η_{sphe} that are conservative, as double-photoelectron emission would result in a very modest increase. Figure 6 shows an example distribution of sphe POD heights that has $\eta_{\text{sphe}} \sim 0.95$. The distribution of η_{sphe}

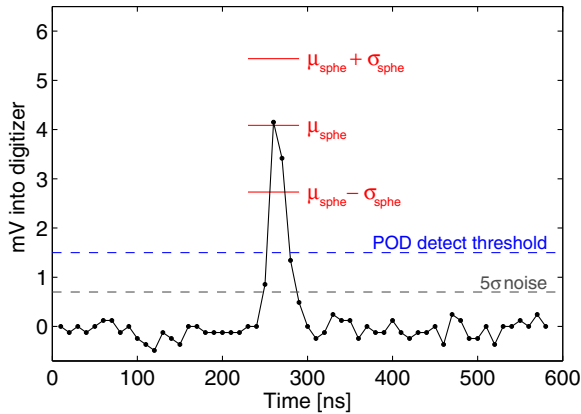


FIG. 5. A schematic diagram of the various thresholds applied to raw waveforms.

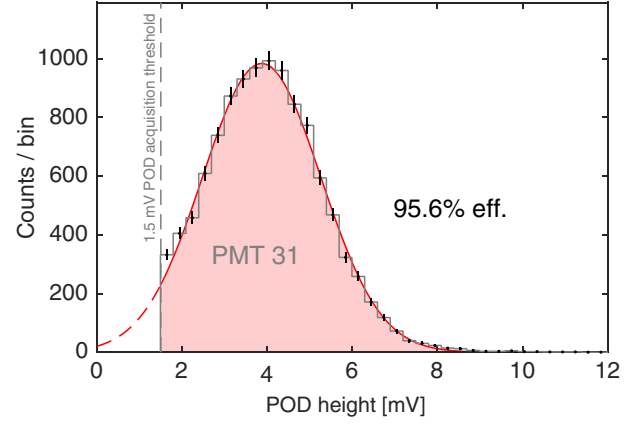


FIG. 6. Measured maximum sphe pulse height for PMT 31. The 1.5 mV POD threshold is indicated by the vertical dashed line. Based on a fit of the single-photoelectron pulse height distribution, it is estimated that 95.6% of the sphe in PMT 31 produce a pulse that exceeds the POD threshold.

values for the PMTs used during Run 3 is shown in Fig. 7. All but two PMTs had a sphe digitization acceptance greater than 0.90; the mode of the distribution is above 0.95.

B. Trigger

The LUX trigger system is described in detail in Ref. [31]. It is a digital field-programmable gate array (FPGA)-based system that flags events in the DAQ data stream for further analysis. All 122 PMTs are summed into 16 trigger channels, with no adjacent PMTs belonging to the same sum. The sums are individually processed using two eight-channel DDC-8DSP digitizers/processors that communicate with a common trigger builder to make a final trigger decision. Internal digital filters perform baseline subtraction and signal

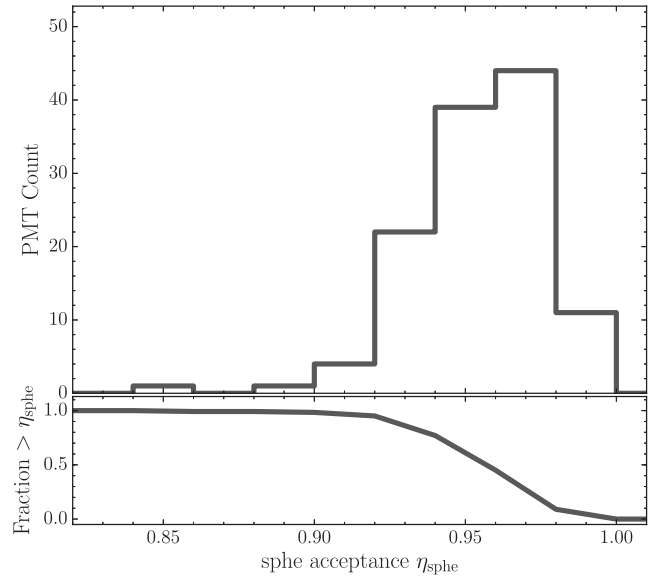


FIG. 7. Single photoelectron acceptance distribution for all 122 photomultiplier tubes.

integration. In Run 3, the trigger required that at least two of these channels have a signal greater than eight photoelectrons within a $2\ \mu\text{s}$ window. The overall trigger efficiency has a strong dependency on the required two-fold coincidence and reaches $> 99\%$ efficiency for S2 signals with a total area of 100 photoelectrons (phe). A detailed study of the trigger efficiency is discussed in Ref. [32]. For the first third of Run 3, the hold-off time, which is a post-trigger time window during which we did not accept new triggers, was set to 4 ms. Analysis of the data being collected showed that the hold-off period did not have to be this long and subsequently was set to 1 ms for the remainder of the run. This improved the live time slightly and was verified to have no negative impact on the overall results.

C. LUX data-processing framework

The LUX data-processing framework (DPF) is a flexible, modular, and multilanguage framework developed by the LUX Collaboration for extracting the relevant features from the raw digitized PMT data and returning a set of reduced quantities that can be used for physics analyses. The LUX DPF employs interchangeable algorithm modules with standardized inputs and outputs that perform predefined tasks, such as calibration, pulse finding, classification, and interaction vertex position reconstruction. These algorithm modules can be written in any supported programming language, which currently includes C++/ROOT, PYTHON, and MATLAB. A MySQL database (referred to as the “LUG”) stores version-controlled calibration values and correction maps, data-processing input settings and data-processing logging, among other bookkeeping values. The LUX DPF was written entirely in PYTHON and can be executed on computing clusters and desktop and laptop computers. The list of modules to use, their order and specific configuration (e.g., threshold value for a pulse-finder algorithm), and required calibration constants must be provided to the framework in an input settings file in XML format—which is stored in the LUG database and associated with a unique identifier (called the global setting number). This identifier allows the collaboration to easily establish the exact conditions that were used to process any data set and ensures that all the data used in a particular analysis campaign has been processed with the same settings.

Figure 8 shows a schematic representation of how raw data are processed through the LUX DPF to obtain the reduced quantities that are used in higher-level analyses. The Event Builder reads the raw data as digitized by the DAQ and extracts the portions that are located in a time window before $\Delta t_{\text{pretrigger}}$ and after $\Delta t_{\text{posttrigger}}$. This window is referred to as an “event.” For Run 3, both $\Delta t_{\text{pretrigger}}$ and $\Delta t_{\text{posttrigger}} = 0.5\ \text{ms}$. Given that the maximum Run 3 electron drift time was $322\ \mu\text{s}$, these values ensure that both the S1 and the S2 pulses are contained in every event since either pulse type may induce a trigger.

The settings used in the Event Builder are stored in the LUG database, and the output file set is assigned a unique

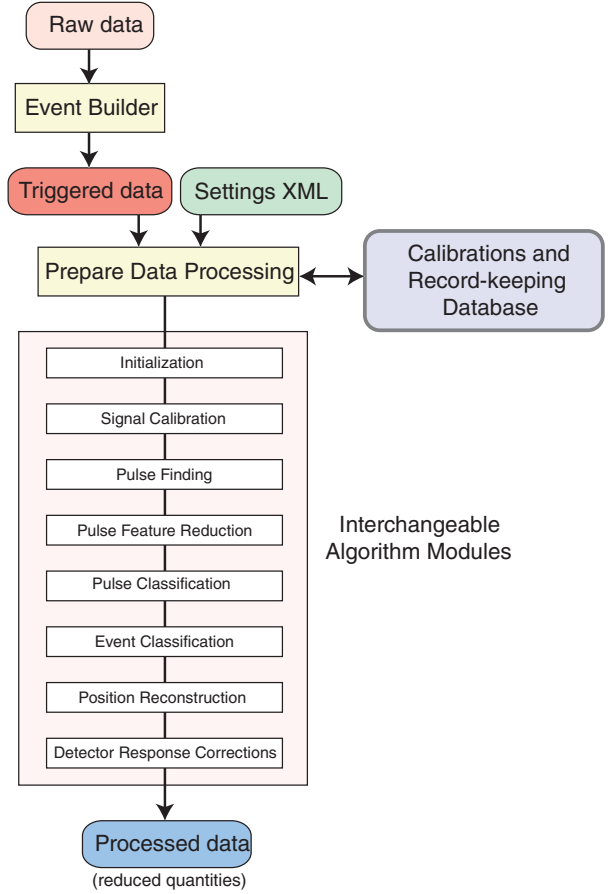


FIG. 8. Schematic view of the LUX Data Processing Framework.

identifier that corresponds with the LUG record. The output of the Event Builder is read by the LUX DPF modules.

The LUG database, in addition to storing data-processing bookkeeping values (such as the Event Builder settings, DPF global settings, and details about each individual data-processing run), also stores calibration constants for the detector. These include PMT gains; x , y , z spatial calibration maps for S1 and S2 pulses; the electron lifetime, detector tilt measurements, light response functions for position reconstruction, and energy calibration parameters, among others. These are sent to each data-processing run as specified in the XML settings file for access by the algorithm modules. The calibration constants stored in the LUG are stored with submission dates, version numbers, originating data set name (from which the values were calculated), and algorithm names. The latter allows for different methods of obtaining a calibration parameter to be selected during a data-processing run.

D. Data-processing algorithms

1. Pulse finder and classifier

At the heart of a reduced-quantities-based analysis, there is a pulse finding algorithm that searches for valid pulses in

the acquired waveforms and which then stores these relevant data for further analysis. Any detected pulses are subsequently classified according to their shape and properties. The pulse finding algorithm module must fulfill the requirements of finding and separating all S1- and S2-type pulses, returning their start and end times accurately. It must do so without excluding a significant fraction of the pulse, as that would lead to under-representation of area. The pulse finder must also respect a limit imposed by the processing conditions of the DP framework of a maximum of 10 stored pulses, a limitation verified through extensive hand-scanning campaigns as not imposing significant efficiency loss.

The pulse finder is based on a sliding boxcar filter (width of $4 \mu\text{s}$), applied to the full event waveform for each triggered event, determining the region that maximizes the enclosed area. A moving average filter (30 ns width) smooths the regions before and after the maximum amplitude in the boxcar. The start and end times of the pulses are set at the point where the smoothed waveforms stay below the set baseline noise threshold of 0.1 detected photoelectrons (phd) per sample for a time of $0.5 \mu\text{s}$. In addition, valid pulses need to be at least 30 ns (three samples) wide, or their average amplitude must exceed the baseline noise threshold. If the pulse width exceeds $6 \mu\text{s}$, a moving average filter with a larger width parameter (250 ns) is used, moving forward and subsequently backward in time from the point of maximum amplitude in the waveform. If a falling edge is followed by a continuous rise in amplitude for a minimum of $0.5 \mu\text{s}$, this filter allows identification of possible additional signal clustered with the original pulse. If two or more individual signals are found, the algorithm splits the original waveform and stores only the start and end time of the largest of the two pulses. Finally, if a pulse has been found, then the corresponding amplitude in the waveform is set to zero, and all of the steps described above are repeated until all pulses in the waveform, up to a maximum of 10, have been identified.

The possibility of missing small S1 signals in the presence of single electrons, PMT afterpulsing, or other spurious signals is minimized by using a complementary search logic that takes advantage of the preponderance for such pulses to dominate the region of the waveform following a large S2 signal. After the first two largest pulses have been found, the waveform is split into two search regions of differing priorities. The region before the pulse that occurred at the later time in the waveform is scanned first with the standard boxcar filter algorithm. If the maximum allowed number of pulses has not been reached after scanning this first search region, then the algorithm will continue to fill the empty slots with pulses from the second search region. This methodology assists identification of events in which there are multiple interactions, i.e., multiple scatters. Figure 9 shows an example waveform of a multiple-scatter event, indicating the regions of the search logic. Once pulses have been identified,

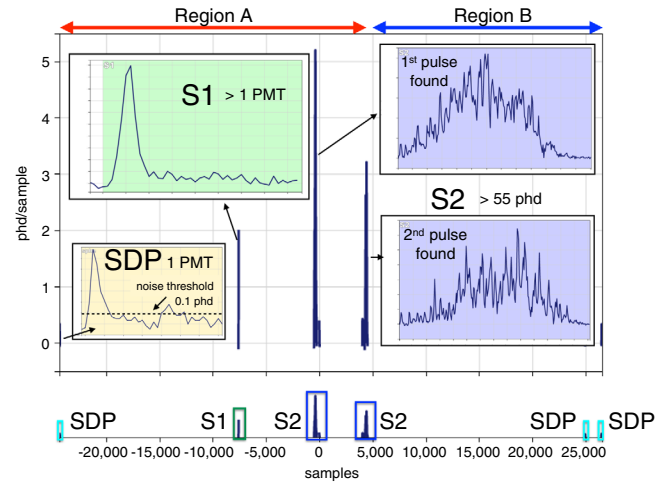


FIG. 9. An example of a double-scatter event waveform. The search regions are labeled A and B, with the first, A, prioritized. In addition to the main signal, this event contains three SDPs, with one shown as an inset to indicate the noise threshold of the pulse finder. The bottom panel shows the location of pulses within the entire event record, with the x axis as the number of 10 ns data samples either side of the trigger pulse, located at zero.

independent modules subsequently parametrize the pulses for further analysis and classification of signal types.

The pulse classification module assigns each identified pulse to one of the following five signal types, according to the extracted pulse parameters: S1, S2, single detected photon (SDP), single electron, and unclassified. The algorithm is represented by the decision tree diagram in Fig. 10. First, S2- and single-electronlike signal types are assigned, followed by S1- and SDP-type signals, ensuring that S1-like signals are not overwritten by S2-like ones. Events that fail all four previous categories are assigned an “unclassified” flag. The signal-type assignment is based on a number of pulse parameters as follows: a width selection, exploiting the much narrower width of an S1- in comparison to an S2-type signal, utilizing the ratio between two boxcar filters of length 0.5 and $2 \mu\text{s}$; pulse rise time, based on width ratios and area fractions, using the distinct shape of an S1-like signal from scintillation

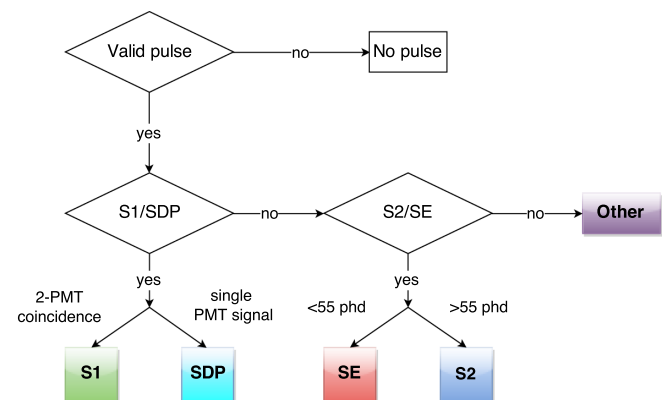


FIG. 10. Pulse classification decision tree.

that features a sharp rise at the pulse start in comparison to the rather symmetrical S2-like signals from electroluminescence in the gas; and PMT hit distribution, where S1 signals are predominately recorded by the bottom PMT array due to total internal reflections off of the liquid-gas interface and the geometry of the TPC. The separation of SDPs and S1s, and likewise single-electrons (SEs) and S2s, is an analysis classification dependent on signal size (rather than being physics based) that enhances coincidental background rejection. For the assignment of an S1 signal, two additional requirements have to be met. First, to reject pulses that are composed solely of baseline fluctuations, the maximum amplitude per sample within the full length of a pulse for at least two individual PMTs must exceed a 0.09 phd/sample amplitude threshold. Second, a two-PMT coincidence requirement is imposed to reject designation of single detected photons from spontaneous PMT photocathode emission as S1 signal. The two-PMT coincidence interval is set to ten samples in width since it is expected that $\sim 80\%$ of the S1 signal for over 90% of both ER and NR pulses will arrive within 100 ns. The two-PMT coincidence requirement must be satisfied by at least two nonadjacent PMTs, each exceeding a 0.3 phd integrated area threshold. A further discrimination between single-electron-type and S2-type signals is gained by recognizing their separation in energy, such that a valid S2 signal must exceed 33 phd in size (corresponding to ~ 1.5 electrons) and classifying those pulses that fall below this threshold as single electrons (as long as their area is greater than 5 phd). Note that further (analysis-dependent) thresholds are usually applied on the S2 size during event classification and selection.

2. Event classification

The event type of interest for the WIMP search analysis is a single-scatter, or “golden,” event. The definition and procedure for the selection of golden events is as follows:

- (i) There is only one valid S2 signal in the event after a valid S1 (golden events may include additional S2s that occur before the selected S1).
- (ii) For the S2 to be valid, it must contain more than 55 “spikes” as counted by digital spike counting (see Sec. III D 4).
- (iii) There is only one valid S1 signal before the selected S2 (S1s following the S2 signal are allowed).
- (iv) The area of the S2 pulse must be larger than the area of the S1.

The purity of the golden events selected by the DPF, and the efficiency of the algorithms used, were evaluated through a detailed hand-scanning campaign. From an americium-beryllium (AmBe) NR calibration data set (live time of 2.53 h), 4000 preselected events were categorized by eye using only the raw waveforms without any information from the reduced quantities. The preselection of events ($\sim 2\%$ of the data set) was necessary to reduce the number to be scanned and, at the same time, to increase the

sample size for single-scatter events in the region of interest. The criteria for preselection were predominately based on event information (number of nonempty samples in an event and the full event area), utilizing only very basic additional pulse parameters such as the largest pulse found and whether a clear subcathode event had been identified. Of the 4000 preselected events, approximately 200 single-scatter events were identified. These were then compared to the result from the event classification DPF module. Applying all WIMP search analysis cuts (see Sec. VI A), the purity of genuine golden events selected by the DPF was determined to be 98%, and the efficiency of the DPF to select golden events from those identified in the data by the hand scan was 98.8%.

3. Lateral position reconstruction

The xy position of an interaction in the LUX detector is recovered directly from the observed S2 signal, by considering the distribution of pulse areas in the upper photomultiplier array. The algorithm used is called “Mercury” and is based on the method developed for the ZEPLIN-III experiment [33]. The algorithm is a statistical search for the xy position that matches the distribution of observed S2 pulse areas in the PMT array to those obtained using a predetermined set of empirical functions, called light response functions (LRFs), which describe the average response of each individual PMT as a function of the interaction position. The major advantage of the Mercury method is that it needs only measured data, rather than simulations, to recover the position of interactions, and thus it can recover features from the data that are not well simulated.

The LRFs for each PMT are obtained through an iterative fit to experimental data, which in the case of Run 3 corresponds to calibration data obtained after the injection of $^{83\text{m}}\text{Kr}$ to the detector (see Sec. V). In each iteration, new LRFs are obtained by fitting the response of the individual PMTs as functions of the event positions. These new functions are then input to the position reconstruction program to derive improved estimates of the position of interactions, which are then used to find new improved LRFs. This process is iterated several times until the functions are stable. For the first fitting iteration only, the initial xy positions are obtained using a simpler method of position reconstruction (e.g., a weighted mean, as used here).

The simplest model for the LRFs, \mathcal{H}_i , consists of a radially symmetric function, η , that depends only on the distance between the center of the PMT and the position of emission of light, ρ . It can be described as

$$\mathcal{H}_i = \mathcal{A}_T \mathcal{C}_i \eta(\rho), \quad (1)$$

where \mathcal{A}_T is the total pulse area of the event and the coefficients \mathcal{C}_i normalize the response of each PMT in the top array [34]. While this model was successfully used in

ZEPLIN-III, it provides unsatisfactory results in the present work, due to the higher reflectivity of the inner walls. In LUX, the inner walls are covered with PTFE, which is a very good reflector for the xenon scintillation light, causing the amount of reflected light to increase for events near the walls. Consequently, a more sophisticated model that used two-dimensional functions for the PMT response has been employed, which models the LRFs as sums of a radial component and a polar component, ε_i , defined for each PMT as

$$\mathcal{H}_i = \mathcal{A}_T \mathcal{C}_i [\eta(\rho) + \varepsilon_i(r, \rho)]. \quad (2)$$

The first component describes the light that propagates directly from the interaction to the PMT and depends only on ρ . The second component corresponds to the light that is reflected in the walls of the detector and is described as a function of the event radial position and also the distance between the event and the center of the PMT [34].

Figure 11 shows the final xy distribution of events in the krypton calibration data set using the Mercury algorithm. The observed striped pattern is a consequence of the geometry of the wires of the grid used to provide the electric field. The pattern originates from a focusing effect of the electrons near the gate wire region, caused by the difference between the fields in the drift region (180 ± 1 V/cm, systematic uncertainty on the mean) and in the liquid extraction region (2.84 ± 0.16 kV/cm).

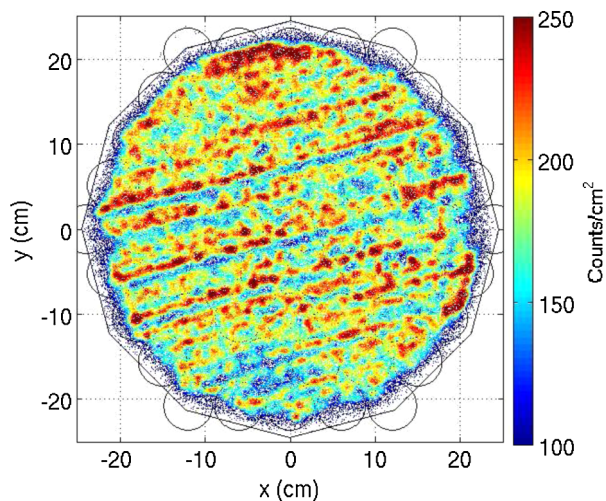


FIG. 11. Map of reconstructed positions in the xy -plane from data obtained after an injection of $^{83\text{m}}\text{Kr}$ in the detector. The observed striped pattern is the result of a focusing effect of the electrons near the gate wires, created by the difference between the fields in the drift and the extraction regions. There are significant differences in the intensity of these strips. These variations are most likely to be caused by small variations in the pitch of the wires producing fluctuations of the bulk field which are not observed for the S2 right below the gate wires [34].

This effect is a limiting factor for the position resolution and may be used to assess the quality of the reconstruction. As discussed later, a nonuniform field within the xenon volume results in events at greater depth having their positions reconstructed at less than their true radius, and it might cause regions with a higher concentration of events; the correction of this effect is discussed separately in Sec. VC.

4. Estimators of light and charge

The S1 signals generated by WIMP recoils are expected to have a mean number of detected photons per PMT that is significantly less than 1. This small signal may be reconstructed with improved precision, as compared to the simple pulse area, by counting of the candidate single-photon pulses in individual PMT waveforms, termed here “spikes.” Such digital counting avoids the variance that is inherent in the width of PMTs’ single-detected-photon area distributions. The raw number of spikes is obtained in LUX by counting maxima in the regions of waveform that are above the POD-start threshold. A simple Monte Carlo model of spike overlap in time, based on average S1 and single-photon pulse shapes, was used to generate a look-up table of the most likely number of true photons as a function of raw area and as a function of counted spikes. This look-up table is then used to estimate the true number of photons comprising the S1, thus correcting for the possibility that photons may overlap in time in a single channel and therefore produce only a single spike. In order to account for the depth of the event, this correction is done separately for the top and bottom arrays. Data from tritiated methane calibrations (see Sec. V) were used to demonstrate that the average of the simple area estimator to this combined area estimator agreed systematically to within 5% everywhere and to within 1% from 16 to 80 phd, above which areas alone were used as pileup became significant.

The last step of event reconstruction was to account for spatial and temporal variation in the detector response. The dominant sources of nonuniformity were the geometry-dependent probability for signal photons to be detected by the PMTs and, additionally for S2, the time-varying concentration of impurities in the xenon which captured ionized charge as it drifted toward the liquid surface. Using the krypton calibration data, maps of the relative position-dependent response were generated for all events. New calibrated S1 and S2 signal areas were defined such that they would equal the raw pulse area for these events if they had occurred at the detector center and there had been no signal loss due to impurities. The symbols $S1_c$ and $S2_c$ were used to denote these final, flat-fielded estimators of light and charge; S2 was always measured via pulse areas, with S1 also utilizing digital spike counting for small pulses (up to 80 phd).

The absolute energy scales of scintillation photons and ionization electrons were obtained from a set of responses to monoenergetic ER sources using the Platzmann model (see Sec. [VD 1](#)); however, the WIMP search did not rely on these scales as the detector's NR and ER responses in $S1_c$ and $S2_c$ were calibrated *in situ*.

Collectively, the above procedures resulted in high-level output of the data-processing framework that was a set of observables measuring position, light, and charge for each reconstructed event above threshold in the active region of LUX: x , y , z , $S1_c$, and $S2_c$. Subsequent analyses applied appropriate event selection and made inferences about physics models by comparing observed and predicted distributions in these observables.

IV. SIMULATIONS: LUXSIM

A. Infrastructure

A detailed understanding of the physics capability of advanced instrumentation is frequently achieved through use of a sophisticated Monte Carlo computer simulation of the apparatus. This is useful in both design and exploitation. Here, the LUX simulation package is presented, known as LUXSIM [\[35\]](#). Its structure can be divided into five overarching, mostly serial functions:

- (i) Recording of simulated energy deposits. This may be for several different particles in multiple volumes at different times and uses the GEANT4 toolkit [\[36,37\]](#), specifically GEANT version 4.9.4 patch 4.
- (ii) The production of both Vacuum Ultra-Violet (VUV) scintillation photons and thermal ionization electrons. This is modeled with the Noble Element Simulation Technique (NEST) [\[38,39\]](#) formalism, a frequently updated semiempirical collection of models based on past detectors' calibration data. The number of such photons and electrons that are generated is in general anticorrelated and depends on the interacting particle type, the magnitude of the electric drift field, and the energies of the recoils.
- (iii) Propagation of photons using the GEANT4 optical model (default libraries within the version specified above). Photons from the initial primary scintillation in the liquid are propagated until reaching PMTs or becoming absorbed by impurities. This is simulated directly by means of either an exponential mean free path for photon propagation through the xenon or by imperfectly reflective surfaces.
- (iv) Drifting of ionization electrons. Low-energy electrons liberated by an interaction are drifted up through the xenon using NEST, diffusing in three dimensions and being absorbed in a manner similar to the photons, but now using the empirically determined (with ^{83m}Kr) electron absorption length, and the field-dependent, binomial electron extraction efficiency at the liquid-gas interface. Once the electrons reach the gas, then the NEST produces the secondary scintillation as a function of field, density, and gas region length. The electroluminescence photons are again simply propagated with GEANT4 optical processes. The drift and extraction fields are both sufficiently uniform in LUX that modeling them as scalar constants serves as an accurate representation for most purposes. See Sec. [VC 3](#).
- (v) Pseudodata generation. Quantum efficiencies (QEs) are simulated as a function of incoming photon wavelength, angle, temperature, and PMT variation. A full custom simulation for the unique DAQ takes numbers and arrival times of primary and secondary photons and generates output files from LUXSIM that are in an identical format to empirical data. These may then be processed using the same data-processing framework, enabling direct comparisons.

LUXSIM provides a component-centric approach; this makes it possible to define any parts of the detector, not just the PMT arrays, as sensitive volumes, making testing and validation studies easier to perform. Although the NEST framework, based on earlier experiments' results, is used for the absolute photon and electron yields, slight modifications are made to the values of various free parameters, fine tuning them to more closely match LUX calibration data at its particular electric field, for both ER and NR, as discussed in Sec. [V](#).

Full ER and NR simulations were performed and passed through the data-processing chain. Initially, the simulation was validated in terms of the ability to reproduce raw waveform shapes, followed by the reproduction of calibration results. With this achieved, LUXSIM was then used to produce ER and NR energy spectra for noncalibration scenarios. Examples included generating samples of "pure" single-scatter events, NR events with no contribution from ER contamination, misidentified multiple-scatter fiducial events, or multiple scatters with vertices outside of the drift region. These are all populations of events that can occur, especially with AmBe or ^{252}Cf sources, demonstrating that the simulation allows exploration of specific (and/or rare) event topologies. The simulation was then also used to provide estimates of the expected WIMP search spectra overlaid with the dominant background contributions from (measured) radioactive impurities in the detector components, contributing ER-type signals overwhelmingly. Predicted WIMP NR signal models could then also be generated, for various candidate WIMP masses and, together with the predicted background spectra, used as probability density functions (PDFs) for use in the profile likelihood ratio (PLR) analysis; see Sec. [VIE](#). LUXSIM combined with the ^{83m}Kr calibration of the position-dependent light collection is critical in simulating WIMP interactions; none of the NR calibrations was uniformly distributed, so they were not direct WIMP analogs.

Extensive campaigns of radioactivity screening with high-purity germanium detectors, inductively coupled plasma mass spectrometry, and neutron activation provided detailed knowledge of the radioactive contamination of many of the detector components, crucial to the high fidelity of any simulation. The results of these studies formed the LUX background model [40] and provided input into the Monte Carlo in terms of radioisotope levels in each component. Simulated S1 and S2 spectra from these contributions were then compared with the actual background observed during the WIMP search, in terms of not only the absolute count rate but also the position-dependent profile, crucial for the PLR analysis (Fig. 12). The primary background constituents are ERs from the PMTs' uranium and thorium decay-chain gamma rays, with further contributions from internal sources (such as ^{127}Xe) and ^{40}K . To precisely match data, the initial contamination values required fine-tuning, though well within errors, from the predictions based upon material screening values.

The simulated sphe area (in voltage) and shape (roughly Gaussian) were calibrated to sphe events occurring in LED calibration and WIMP search data sets. In principle, the size of single electrons, in terms of number of phe, could have been predicted by calculating the absolute S2 photon yield [41,42] and knowing the light collection efficiency for the gas region from optical simulations and the measured QEs. This method in fact agrees within 10% of the observed single-electron pulse area [43,44]. The width and general shape are matched using a Gaussian distribution for the absolute yield in the gas gap and binomial light collection and QE, yielding a slightly non-Gaussian shape that is indeed observed in the data. A comparison of the single-electron area between simulations and data is shown in Fig. 18 and discussed further in Sec. V. The single-electron width in the phe number is about twice the value expected from a Poisson process (\sqrt{N}), though this is understood as being due to field nonuniformities between the anode and liquid-gas border [45]. The temporal profile of a single electron, or of an S2 of any size in general (see Fig. 13), matched well that expected given the known electron diffusion constants in liquid xenon measured in previous experiments decades ago (extrapolated to low field from Fig. 12 of Ref. [46]), the electron drift speeds in liquid and gas, the electron extraction delay at the liquid surface, the (small) light travel time to the PMTs, and the singlet and triplet time constants characteristic of the excited molecular (excimer) states of xenon dimers. The last two contributions, plus the time it takes for ionization electrons to recombine with ions, also allows reproduction of the S1 pulse shape [47,48].

B. Optical model

In contrast to NEST, the model for photon propagation has been specifically tailored to the present detector conditions. The parameters within the GEANT4 model that

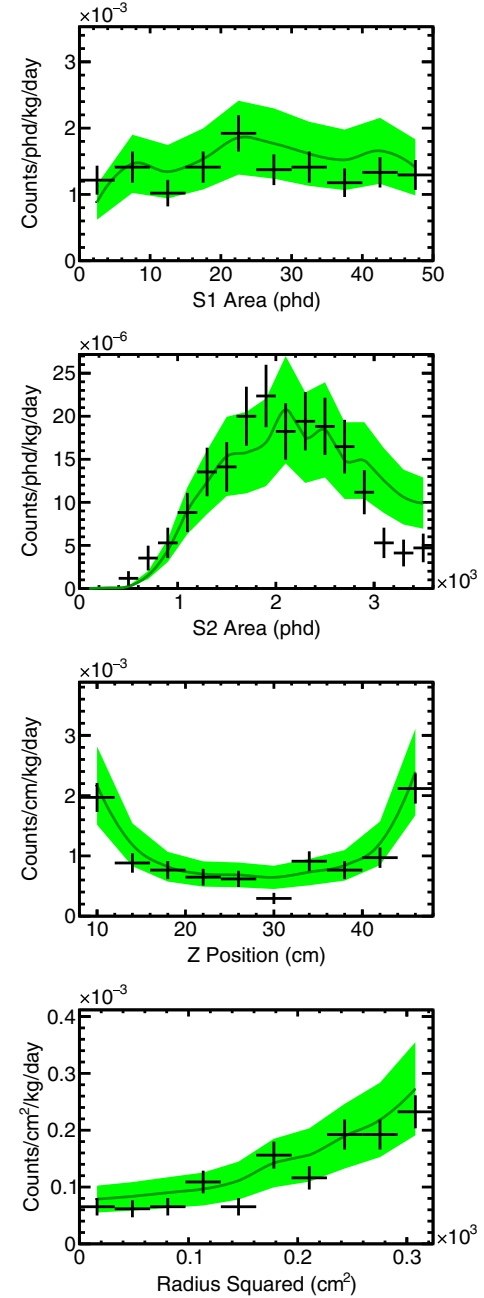


FIG. 12. From top to bottom: Profile of the absolute number of background events as a function of S1 size, S2 size, depth, and radius, combining all known sources of background. S1 is cut off at 50 phd (photons detected), the WIMP search cut, causing a dropoff in S2 spectrum. The black points and crosses are the data, while the smooth gray center lines come from a simulation based on a data-informed full background model. The green bands represent the 1-sigma uncertainties in the simulation; they are $\sim 30\%$ of the value of the background model, aggregating uncertainties across different background components [40]. The disagreement at high S2 comes from the NEST overestimating the high-energy charge yield for electron recoils. Because this was at high energy, the impact on the WIMP search sensitivity was negligible. (This was corrected in the following run, 332 live days, with new calibrations.)

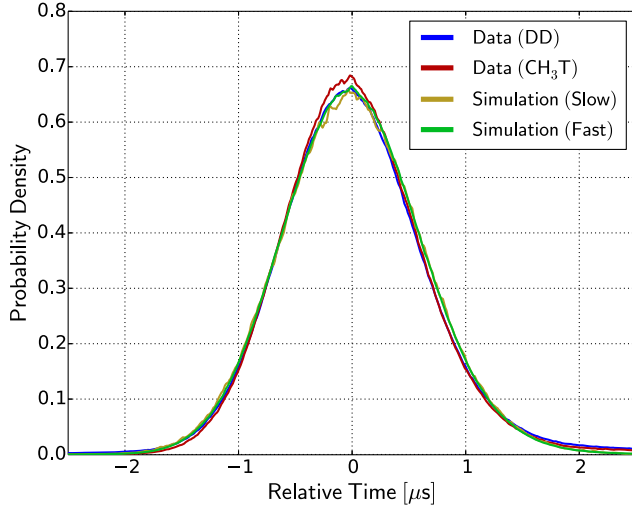


FIG. 13. The averaged S2 shapes from NR (D-D neutron calibration data, in blue) and ER (CH_3T , in red). Slow simulation (yellow) refers to a G4/NEST prediction from first principles based on the electrical and geometric properties of the detector. Fast simulation (green) refers to the parametrized PMT hit-map library mode of LUXSIM that is forced to agree with data as closely as possible. A drift time cut of 100–120 μs is applied in all cases so that D-D and tritium can be compared directly. In contrast to S1 signals, no significant differences are expected for different interaction types.

were tuned, such that the simulation accurately replicated the data were:

- (i) the bihemispherical reflectance of the stainless steel field-generating grids assumed to be constant with the angle of incidence and the bihemispherical reflectance of PTFE in liquid and in gas assumed to be only diffuse (Lambertian).
- (ii) the photon mean free path, separately tuned for liquid and gas phases.
- (iii) the reflectivity of the aluminum flashing on the PMT quartz faces.
- (iv) the average QE for the top PMT array relative to the bottom, an absolute normalization systematic used to provide better agreement with data for the ratio of top/bottom signal size. This correction, around 2%, was within the expected uncertainty.

Table I shows the optical model parameters used in LUXSIM optimized for replication of the experimental data. The indices of refraction of xenon and quartz were based on previous data [49,50], while the Rayleigh scattering length came from a theoretical calculation [51]. Of these, the parameters with the most impact were found to be the reflectivity of the PTFE walls in the liquid and the photon absorption length in the liquid. For both, an increase in the parameter value implies better light collection, and thus they are partially degenerate, although some discrimination was afforded through comparisons of S1, S2, the position dependence of each, and the ratio of light in the top or bottom PMTs to the total light collected (Figs. 15 and 16).

TABLE I. Optical model parameters used in LUXSIM optimized for replication of the experimental data. The predicted QE refers to measurements of LUX PMTs warm, corrected with earlier measurements of the same PMTs indicating an increase when cold.

Quantity	Liquid	Gas
PTFE diffuse reflectivity (%)	97^{+3}_{-2}	75^{+10}_{-5}
Stainless steel grid reflectivity (%)	5 ± 5	20 ± 5
PMT aluminum reflectivity ^a (%)	100^{+0}_{-10}	100^{+0}_{-10}
Photon absorption (m)	30^{+40}_{-20}	6 ± 3
PMT array QE/predicted	1.024	1.000

^aThe aluminum is in contact with the PMT quartz window.

The strongest optimization of the optical model was found to come from consideration of the S1 pulse area as a function of depth from the main monoenergetic peak (32.1 keV) in high-statistics $^{83\text{m}}\text{Kr}$ calibration data. The radial and angular dependencies, due to the orientation of the grids and the variation of PMT QEs, were found to be of secondary importance. The vertical symmetry is significantly broken by total internal reflection at the liquid surface.

After the relative light collection efficiency, as a function of depth and radius, had been matched for the S1 signals with LUXSIM, it was validated through analysis of single-electron signal size and position dependence of S2 signals. The use of a $^{83\text{m}}\text{Kr}$ calibration source led to a model that is valid for all particles because, once light is generated, the nature of the original incident particle is irrelevant. The absolute photon detection efficiency, which is a combination of light collection and QE, was then simulated for the events that occurred at the center of the detector, and this was then used to surpass the relative efficiency comparison. The resulting simulated S1 photon detection efficiency, g_1 (as defined in Sec. V D as the ratio of photons detected to all those emitted), of 0.119 ± 0.001 phd/photon is in excellent agreement with the value of 0.117 ± 0.003 phd/photon determined from analysis of the data presented in Fig. 33. The simulation result of 0.119 phd/photon includes a factor of 1.122 arising from the increase in QE occurring at cryogenic temperatures [52,53]. This is consistent with the previously reported result of $14 \pm 1\%$ [18], which was also based on simulation, but which did not include the recently measured average 17.3% probability for the R8778 PMTs to generate 2 phe from a single detected photon [19,54]. Lastly, comparison of monoenergetic S1 peak mean positions to NEST predictions agreed within known uncertainties (Fig. 14 is one example). The revised calculation of the extraction efficiency (49% compared to the previous value of 65%) is now in better agreement with Refs. [55,56], but differs from Ref. [46], thus leading to a lower extraction electric field than originally estimated.

Deviations between data and simulation in Fig. 15 demonstrate that a purely physics-motivated approach, even one which includes tuned optical properties such as reflectivity, is still imprecise. It cannot, for instance, account for

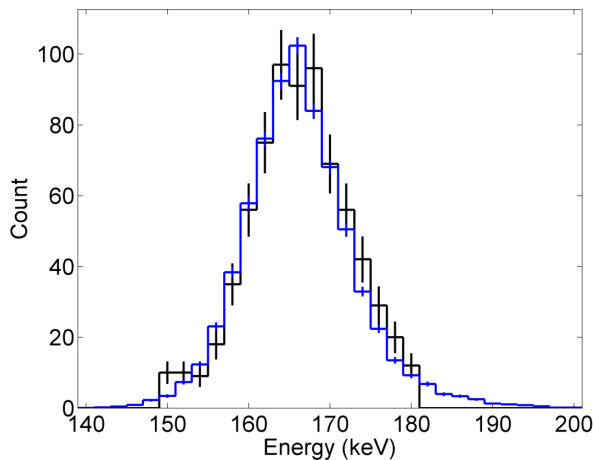


FIG. 14. Comparison of simulation results (blue) to data (black, with error bars representing the statistical uncertainties in each bin) for the $^{131\text{m}}\text{Xe}$ cosmogenic-neutron activation peak, which is the only monoenergetic ER in the background that has sufficient statistics to allow such a comparison to be made. Presented here is the reconstructed “combined” energy scale, calculated from both S1 and S2 and which therefore has improved resolution as compared to the use of a single channel alone, for the data, while for the simulation the x axis represents the Monte Carlo truth energy. The width of the peak is determined by the binomial statistics of S1 light collection, sphe resolution, electron lifetime, extraction efficiency, S2 collection, single-electron pulse size variance, and recombination fluctuations. To reproduce the mean precisely, the values of S1 and S2 collection efficiencies had to be varied by 5% (upward), well within their uncertainties at the 1–2 sigma level, and the recombination fluctuations are increased by 1-sigma compared to a best fit at low energies. The default NEST mean yields from V0.98, 2013, are assumed without modification. The source of the slight asymmetry in the data was likely a background component.

all the potential microscopic surface deformations that create position-dependent reflectivities on a surface or possible exotic nonexponential mean free paths for photon absorption by impurities in the xenon. Thus, to ensure the most precise possible background and signal models for the sensitivity calculation, the position corrections based on the $^{83\text{m}}\text{Kr}$ calibrations (see Secs. III D 4 and V and the data within Fig. 15) applied to the real data sets were also used to generate empirical S1 and S2 three-dimensional look-up libraries for every PMT. These libraries avoided the need for photon propagation in every simulation and thus allowed 2 to 3 orders of magnitude larger simulated data sets to be generated.

In summary, the LUXSIM simulation described here provides near-perfect replication of numerous empirical results, providing exceedingly reliable support for further data analysis.

V. CALIBRATIONS

A. Responses to single quanta

Energy depositions in experiments such as LUX arise from scintillation and ionization, both of which result in

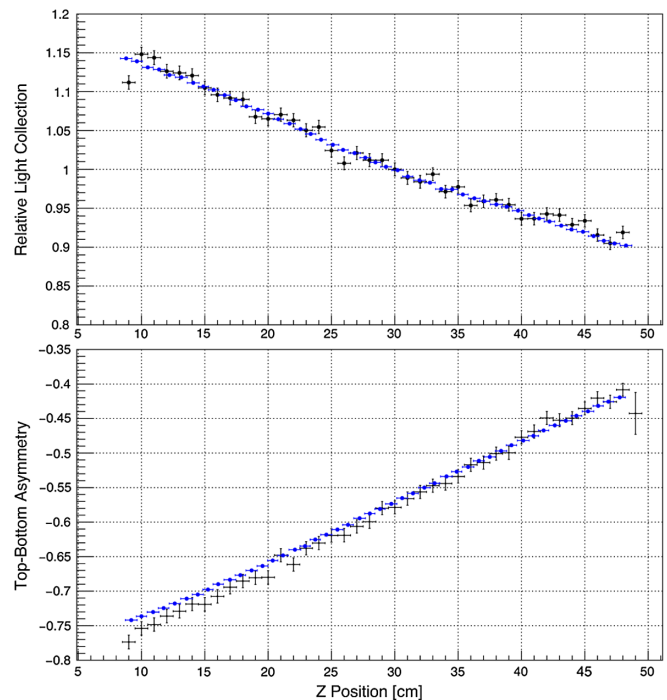


FIG. 15. Top: the relative light collection of S1 signals as a function of the depth in the detector. Bottom: The asymmetry in the S1 light collection efficiencies between the top and bottom PMT arrays, defined as $(T - B)/(T + B)$, where T and B are the values of S1 recorded in the top and bottom PMT arrays, respectively (simultaneous fit to both quantities). In both plots, data are presented in black with statistics-only error bars, while the results of the simulation are shown in blue.

photons detected by PMTs. The variance in such signals is dominated by the number of phe emitted at the photocathodes, as this is where the signal quanta is at a minimum (after the QE and before dynode amplification). Consequently, the basic unit of measurement in LUX is the number of phe. Calibration begins with consideration of sphe. To stimulate the emission of sphe for calibration purposes, six blue LEDs, located in the LXe but outside of the TPC, were individually pulsed at a rate of 1 kHz and a pulse width of 100 ns. The pulse amplitude was set so that a given PMT saw no signal for most LED pulses, in which case the number of phe observed per LED pulse in that PMT obeyed a Poisson distribution. If the amplitude is small enough, LED pulses that show nonzero signal will be a nearly pure sample of sphe and may be used to extract the average amplification (i.e., the “gain”) of the PMT.

The single-phe response of one LUX PMT is shown in Fig. 17 (gray histogram). The gain of the PMT, defined as the average number of electrons collected at the anode from a single electron emitted from the photocathode, is on average $(3.9 \pm 0.4) \times 10^6$, with a resolution (σ/μ) of 35%. LED calibrations were carried out weekly throughout the duration of the experiment.

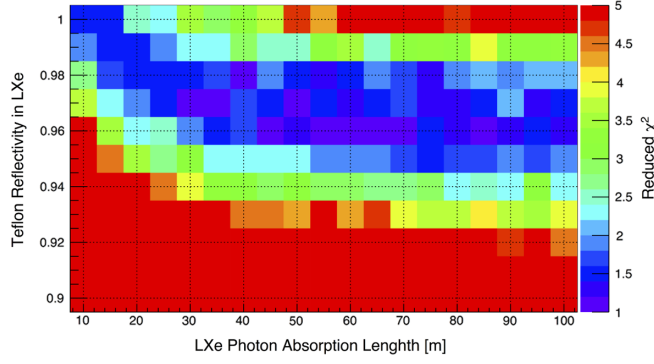


FIG. 16. χ^2 map for Teflon reflectivity and photon absorption length in liquid xenon, the two free parameters that have the greatest impact on both relative and absolute light collection. The degeneracy between the reflectivity of the Teflon and the absorption length for photons in the liquid can be seen, with each having similar effects on the reduction in the light collected. As the former asymptotes to $\sim 96\%$, a reasonable value [57], the latter approaches infinity.

Typically, in the use of PMTs, a sphe is synonymous with a single detected photon. However, as shown in Fig. 17, the LUX experiment has discovered that the detection of a single VUV scintillation photon often causes the emission of two phe. Because the QE reported by the PMT manufacturer is defined as the number of phe emitted from the photocathode per incident photon, this effect leads to a difference between the QE and the photon detection efficiency, a fact that has important implications in accurately reconstructing the energy of an event based on absolute yields [54].

The rate of double-phe emission per VUV photon has been measured in two ways: first using scintillation photons emitted by the liquid and second by using scintillation photons emitted from the gas. For the former, a sample of events from a CH_3T calibration was selected, for which the total S1 light collected for each event is on average ~ 5 phe. With this selection, the average number of phe collected in a single PMT is less than 0.05 per event. The number of detected photons—not the number of phe—is Poisson distributed in this PMT (similar to the strategy utilized in LED calibrations). If μ is the average number of detected photons per event for a given PMT, the fraction of nonzero events which are contaminated by multiple detected photons is $1 - (\mu e^{-\mu}) / (1 - e^{-\mu})$. Therefore, taking $\mu < 0.05$, the set of nonzero hits (for that PMT) is a nearly pure sample of single detected photons, with multiple detected photons contributing less than 2.5%. Figure 17 shows, for an example PMT, a comparison between the spectrum obtained from the (optical) LED calibration and the (VUV) CH_3T calibration [“VUV (liquid)”]. The shoulder on the tail of the single-phe peak is readily visible, which indicates the presence of this double-phe emission process. Plots such as this were used to construct a “VUV gain” for each PMT, which indicated the average number of electrons

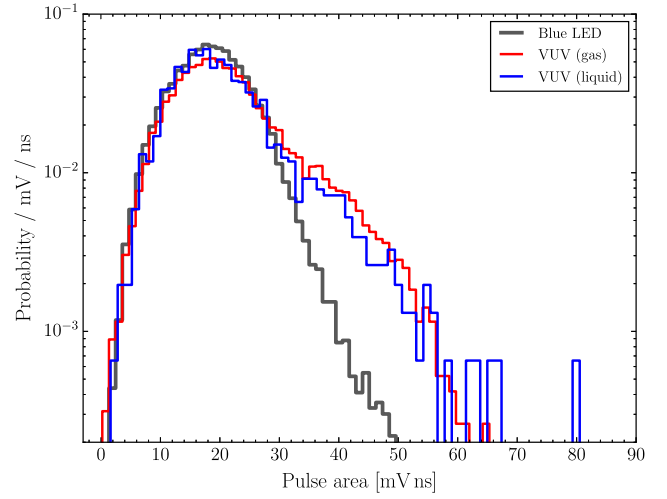


FIG. 17. The pulse-area spectrum for one of the LUX PMTs under a range of conditions. The black histogram shows single photoelectrons stimulated by exposure to a pulsed blue LED, the red histogram shows the signals recorded as a consequence of VUV photons emitted from gaseous xenon, and the blue histogram shows the signals recorded as a consequence of VUV photons emitted from liquid xenon. Evidence of two phe emission from the VUV photon exposures is clearly evident.

collected at the PMT anode for a single detected VUV photon. Consequently, the basic unit of measurement is no longer the number of phe but the number of photons detected, phd.

The second method for measuring VUV-photon response uses electroluminescence light from the gas region, in the form of single-electron ionization pulses from calibration data. Photoionization of impurities in the bulk liquid following $^{83\text{m}}\text{Kr}$ S1s provide a large and pure sample of single electrons. Light from each extracted electron is approximately uniform in time over the $1\ \mu\text{s}$ drift from the liquid surface to the anode and sums to an average of 25 detected photons across the 122-PMT array; the signal therefore appears in individual PMT traces predominantly as single photons or two clearly separable photons (single photons having FWHM around 30 ns). The mean area of the single-photon response in a given PMT is obtained in three steps. First, the mean area of those single-electron responses with one identified maximum above a 1.4 mV threshold is calculated. Second, the number of unresolved pileup events contributing to that mean is estimated from those responses with two photons resolved in time: the interval between first and last DAQ samples above threshold for the two-spike responses has the expected linear distribution above seven samples, which is extrapolated and integrated over the region of smaller intervals, where the two photons may not be resolved. Third, the mean area of single spike events is corrected for the pileup, with the area of the contaminating two-photon responses taken to be the same as the resolved two-photon responses. This correction is small, on average 3% for top array PMTs and 1% for

bottom array PMTs. The resulting gain estimates are systematically 2.5% higher than the liquid-scintillation estimates, which may be due to the difference in scintillation wavelength. The liquid-scintillation values are adopted since it is for S1 light that the number of detected photons implies a detection efficiency for the fundamental signal quanta. It is worth noting that for the case of ionization any pulse-area unit used for both S2s and single electrons cancels out when one divides to estimate the signal size in units of electrons.

In a two-phase Xe TPC, ionization signals generated by a particle (or gamma-ray) interaction result in a burst of VUV photons. The size of these pulses can be understood more usefully by reporting the absolute number of electrons creating the signal. Doing so requires a calibration of the detector to single electrons. Fortunately, single electrons are periodically emitted from the liquid surface, a phenomenon that has been described in the ZEPLIN-II [58], ZEPLIN-III [59], XENON10 [10], and XENON100 [60] experiments. A sample of pure single electrons is selected by searching the event record in ^{83}mKr events between the S1 and S2 signals. Since it is known that these events are essentially single site, any S2 features in the event record between S1 and S2 are likely to be single electrons. The rate of single electrons is low enough that the probability for two electrons to randomly overlap in time is negligible. Figure 18 shows the spectrum of the ionization signal from these single electrons, in units of the number of detected photons. As the proportional scintillation process depends

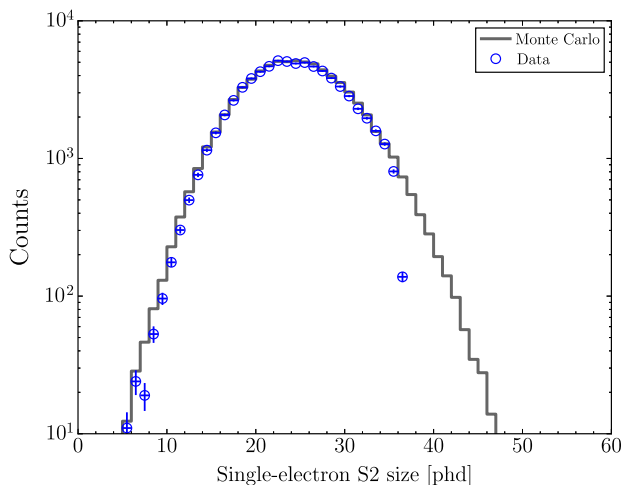


FIG. 18. The spectrum of ionization signals from single electrons, shown as the blue data points. The spectrum, fitted with a skew-normal distribution, gives a mean of $24.66 \pm 0.02(\text{stat})$ phd and a $1-\sigma$ width of $5.95 \pm 0.02(\text{stat})$ phd. The black line shows a simulation of this spectrum produced from first principles, based on a 5.8 kV/cm field in a 1.6 bar, 5 mm “gas gap.” with an S2 light collection efficiency of 10%. The falloff in the data for large phd values is caused by the switch from single-electron to S2 pulse definition.

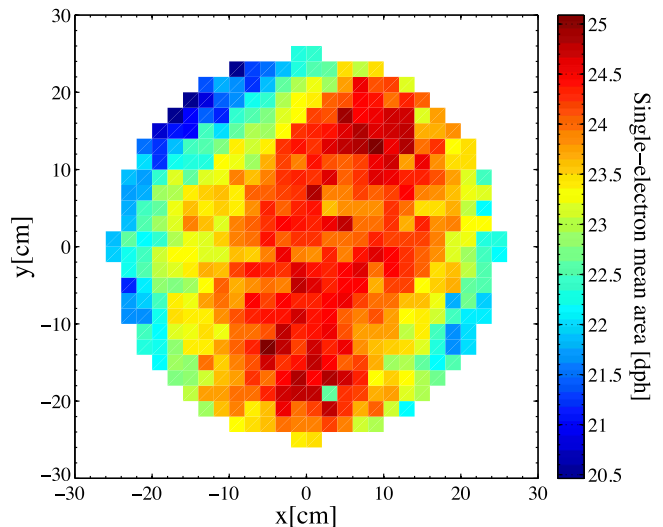


FIG. 19. Variation of the mean of the single-electron distribution in the plane of the liquid surface.

on the extraction field and gas gap [61], variations of these parameters in time and across the plane of the liquid surface can lead to variations of the average single-electron size. Figure 19 shows the average of the single-electron distribution in the x - y -plane, while Fig. 22 shows the average single-electron size over the duration of the science run.

B. Stability

1. PMT gain stability

The LED calibrations, described above, were performed periodically over the course of the WIMP search. Figure 20 shows the deduced gains for ten PMTs over most of the duration of the run, illustrating the stability achieved. For all PMTs, the relative level of fluctuations is presented in Fig. 21; most gains are stable to better than 2% (standard deviation/mean).

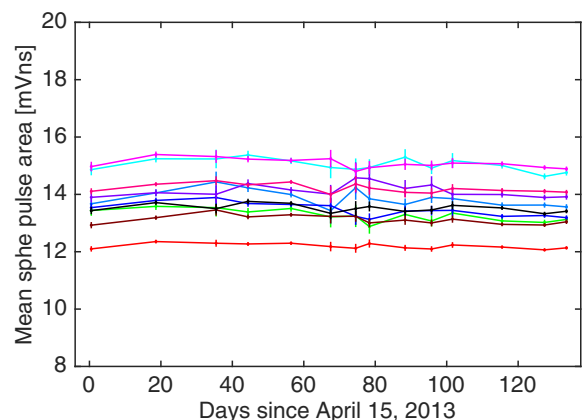


FIG. 20. The gain as a function of time for ten representative PMTs.

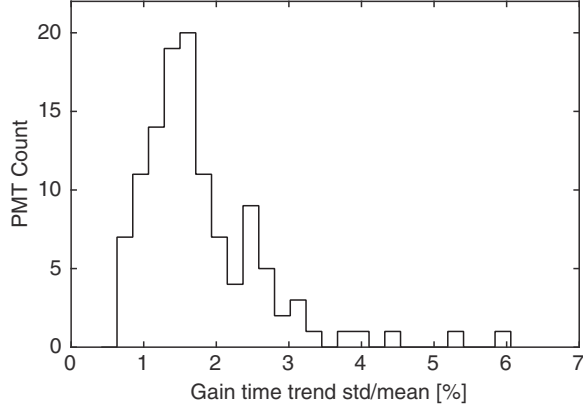


FIG. 21. The gain of each PMT was measured each week throughout the run with an LED calibration to stimulate sphe emission. Shown is a histogram of the standard deviations of these gains, each normalized to the PMT's mean gain. Most gains are stable to better than 2%.

2. Single-electron size stability

The gain for S2 signals, generated by electroluminescence in the gas, is determined by a number of factors. These include the pressure of the gas, the level of the liquid, and the extraction electric field. While each of these parameters was monitored by the slow-control system, an absolute measurement of the S2 gain was determined using $^{83\text{m}}\text{Kr}$ calibration data. As mentioned previously, these data contain single electrons emitted from the liquid surface, allowing the average size of a single electron S2 to be determined periodically over the course of the WIMP search. The resulting gains are shown in Fig. 22. These data indicate that the single-electron size was stable at the level of 1.4% over the duration of the run.

3. Electron lifetime stability

As electrons are drifted through the liquid, they may be captured by residual electronegative impurities, such as O_2 . This produces a depth-dependent attenuation of the ionization signal and allows the concentration of O_2 to be

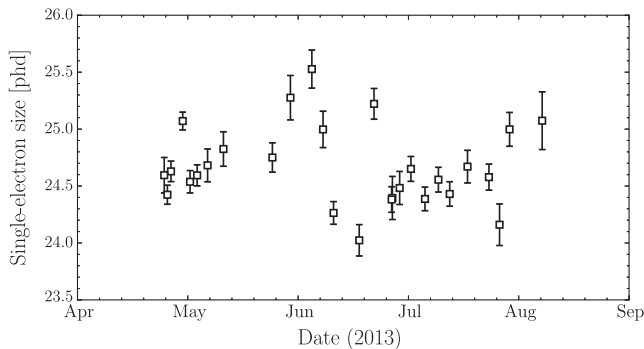


FIG. 22. The variation in the size of single-electron S2 pulses over the duration of the WIMP search.

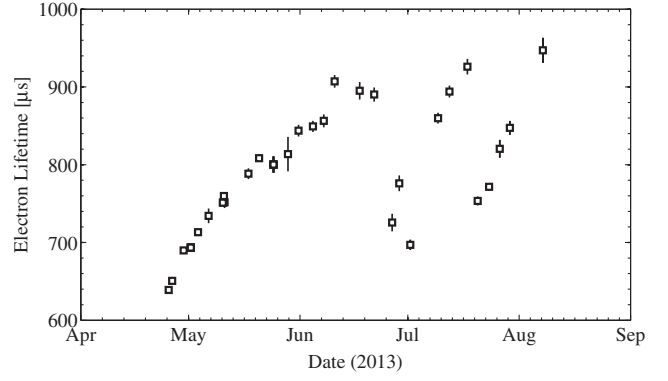


FIG. 23. The measured electron lifetime over the duration of the experiment. Sporadic dips in electron lifetime are a result of temporary stoppages in the LXe recirculation.

expressed as the average time that an electron will propagate before capture; this quantity is known as the electron lifetime. As this quantity varies over the course of the experiment (and is used to correct observed ionization signals), it is measured periodically during the WIMP search. Periodic calibrations with $^{83\text{m}}\text{Kr}$, in a technique similar to those described in Refs. [62,63], were carried out. This source decayed via two sequential transitions depositing 32.1 and 9.4 keV, each of which consisted mostly of conversion and Auger electrons, with small contributions of gamma and x rays. The half-life of the 9.4 keV state was 154 ns, sufficiently short enough that the sequential decays occurred at essentially the same physical location and also that they often occurred too close in time to be resolved as separate pulses. The depth dependence of the combined 41.5 keV ionization signal was measured, from which the electron lifetime was calculated. Figure 23 shows the measured electron lifetime during the WIMP search.

4. $^{83\text{m}}\text{Kr}$ light yield stability

The stability of the scintillation and ionization was monitored with periodic $^{83\text{m}}\text{Kr}$ calibrations [22]. Figure 24 shows these responses over the time period of the WIMP search. The relative time variation (standard deviation/mean) of the scintillation response from these calibrations is 0.6%, while that of the ionization signal is 2.4%.

C. Position reconstruction

1. Measuring the fiducial volume

The total xenon mass in the active volume between the cathode and the gate grids is determined as

$$M_{\text{Xenon}} = 3R^2L\rho_{\text{Xenon}} = 250.9 \pm 2.1 \text{ kg}, \quad (3)$$

where $R = 24.48 \pm 0.05 \text{ cm}$ is the distance between the center of the detector and the corners of the PTFE panels and $L = 48.32 \pm 0.34 \text{ cm}$ is the distance between the

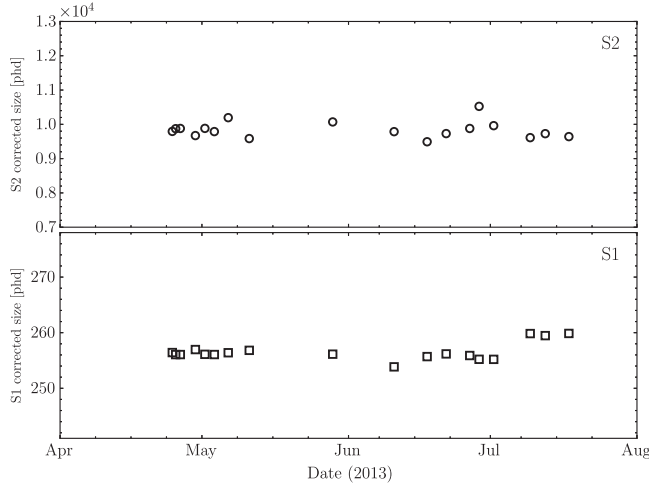


FIG. 24. The scintillation (squares) and ionization (circles) response of the LUX detector over time to the decays of $^{83\text{m}}\text{Kr}$.

cathode and the gate grid (dimensions when the detector is cold). $\rho = 2.887 \pm 0.005$ g/mL is the average xenon density during the run, corresponding to an average temperature of 173.19 ± 0.07 K.

The fiducial region, that is the volume within which candidate events must occur, is defined as a cylinder with radius 20 cm and a vertical extent corresponding to drift times between 38 and 305 μs . The conversion between the drift time and z position, measured relative to the surface of the PMTs in the bottom array, is given by

$$z = s_c + L - v_d(\tau - \tau_g) \quad \text{with} \quad v_d = \frac{L}{(\tau_c - \tau_g)}, \quad (4)$$

where v_d is the drift velocity, measured at 0.1518 ± 0.0011 cm/ μs ; $s_c = 5.6$ cm is the distance between cathode and the surface of the PMTs; and τ_g and τ_c correspond to the drift time of events at the gate and at the cathode, respectively. These values are estimated using the krypton calibration data, from which $\tau_g = 3.4 \pm 0.4$ μs and $\tau_c = 322.3 \pm 0.4$ μs are obtained. The resulting fiducial volume mass is found to be 147 ± 1 kg.

The fiducial mass of the detector may also be calculated using the distribution of the tritium events in the detector. This method takes the ratio of the observed number of tritium events in the fiducial volume to the number of events between the cathode and the gate and then multiplies it by the total xenon mass between the cathode and the gate. This method relies on the tritiated methane being isotropically distributed throughout the detector but has the advantage of not requiring physical dimensions to be defined, such that possible systematic uncertainty from the drift time conversion and position reconstruction are eliminated.

For this calculation, the tritium events are selected using the same quality cuts as applied in the WIMP search data

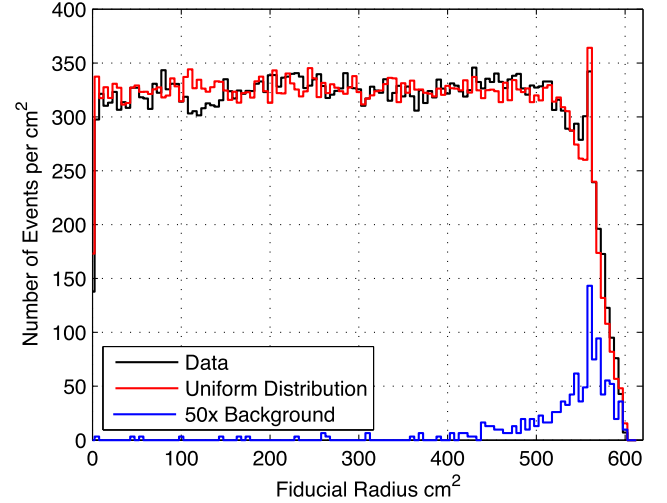


FIG. 25. Histogram of the number of the observed tritium events as a function of the corrected radius and for a drift time between 38 and 305 μs (solid black line). The red line represents a simulation of a perfectly uniform distribution of events. The blue line represents an estimate of the background observed during the lifetime of the tritium calibration, multiplied by a factor of 50.

(described in Sec. VI A). Additionally, the events from the first hour after the tritium injection are not included to ensure uniform distribution of tritium. The mixing time observed for $^{83\text{m}}\text{Kr}$ injections was observed to be less than 10 minutes, and a similar mixing time is expected for tritium injections. After the mixing period, the distribution of tritium events was found to be highly uniform along the radius, with only a small accumulation of events near the walls of the detector, as shown in Fig. 25.

Using a drift time cut to select events in the active volume between the cathode and the gate grid, $(223.1 \pm 0.5) \times 10^3$ tritium events were observed. A fiducial volume cut selected events with a drift time between 38 and 305 μs and a radius less than r . The ratio between these two event counts is multiplied by the xenon mass between the gate and cathode, calculated in Eq. (3). Figure 26 shows the fiducial mass as function of the radius r . For a radius of $r = 20$ cm, the number of events observed is $(129.1 \pm 0.4) \times 10^3$, which yields a fiducial mass of 145 ± 1 kg.

Figure 27 shows the difference between the mass calculated using the event counting method and the mass calculated directly from the geometry, as function of the fiducial volume radius. The difference between these two methods is smaller than 1.4 kg for any fiducial radius.

2. Accuracy of position reconstruction

Given the geometry of the LUX detector, the statistical uncertainty in xy position reconstruction is most naturally handled in polar coordinates. In this case, the uncertainty breaks down into three categories:

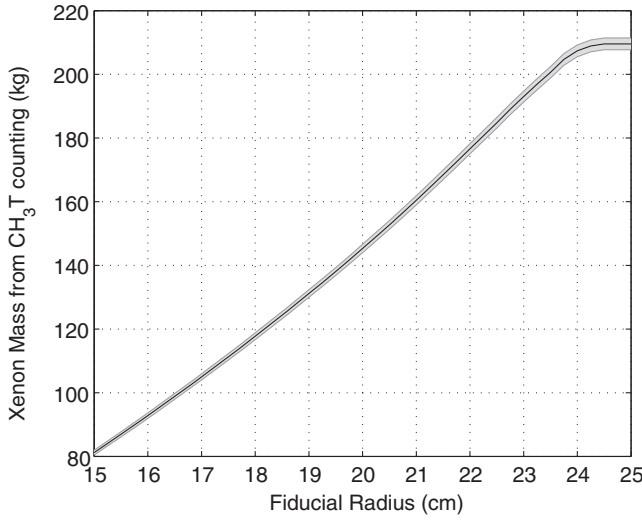


FIG. 26. Xenon mass calculated directly from tritium counting as a function of the radius of the fiducial volume. The ± 1 sigma band is represented by the gray band. The drift time considered is between 38 and 305 μs , and the mass between the gate and anode is 250.9 ± 2.1 kg.

- (i) The centripetal uncertainty: the radial uncertainty in the direction toward zero radius.
- (ii) The centrifugal uncertainty: the radial uncertainty in the direction toward the wall.
- (iii) The polar uncertainty: the uncertainty in the polar angle, treated as symmetric.

This parametrization is convenient as the radial position is paramount in determining whether an event looks more

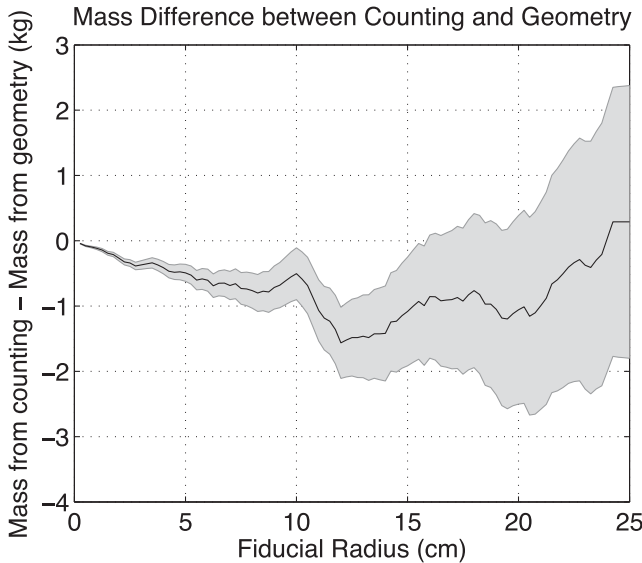


FIG. 27. The difference between the xenon mass calculated from the counting of tritium events and the xenon mass obtained from the geometry of the fiducial volume as function of the radius of the fiducial volume. The $\pm 1\sigma$ region is represented by the gray band.

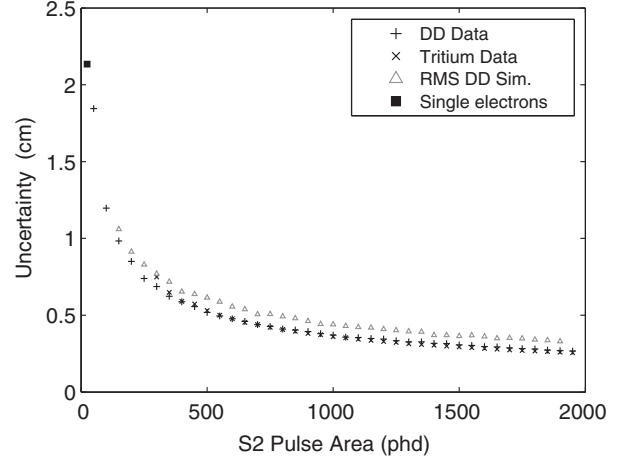


FIG. 28. Comparison of centripetal uncertainties in Mercury in D-D data (vertical black crosses), tritium data (diagonal black crosses), single electrons (black squares), and the root mean square between the reconstructed radius and the original radius, obtained in D-D simulations (triangles).

signal-like (in the bulk, toward the center of the detector) or backgroundlike (toward the wall of the detector). These three uncertainties are calculated from the Mercury error oval based on the goodness of fit. Figure 28 shows the centripetal uncertainty vs S2 size for D-D data and tritium data and the root mean square of the difference between the reconstructed radius and true radius, from D-D simulations. Here, the uncertainties, binned by S2 size, are the average over the entire fiducial volume. Unsurprisingly, the uncertainty decreases as S2 size increases, as a result of there being more information carriers. The uncertainties calculated from the D-D data and tritium data agree between each other, while the results obtained from the D-D simulations are 13% larger. On average, the results for the centrifugal uncertainties are comparable to the centripetal uncertainties.

An advantage of comparing the uncertainty in simulation to that of data is that in a simulation the true event locations are known, allowing for a consistency check, as shown in Fig. 29. The calculated statistical uncertainties are very close to the differences between the real and reconstructed positions.

In addition to the statistical uncertainty in the xy - ($R\theta$) positions, there is a systematic uncertainty that may be calculated using the position of events in the beam from D-D calibrations. Because the D-D events arise from a well-collimated beam, the events in the D-D calibrations appear as a line in the xy position ($R\theta$) plane. The deviation from a straight line may therefore be used as a measure of the systematic uncertainty in the position reconstruction. This measurement of the systematic uncertainty does not depend on the actual angle of the beam, it depends only on the beam being collimated with a small angular divergence. A linear fit to the D-D path and the reconstructed xy

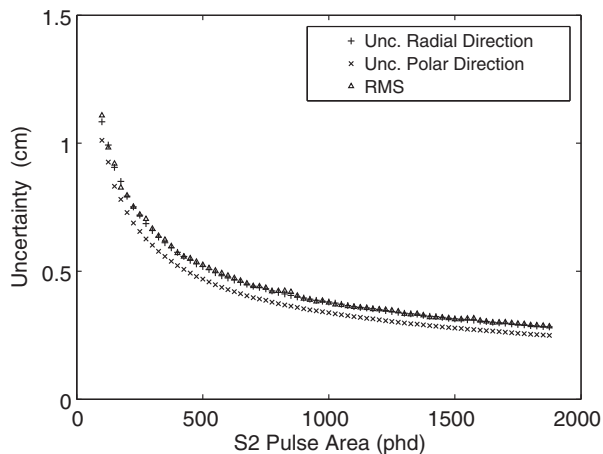


FIG. 29. Comparison of the centripetal and polar uncertainties in D-D simulations to the root mean square of the difference between the reconstructed position and the true position.

position in the D-D data is shown in Fig. 30. The maximum deviation from the linear fit within the fiducial volume is 4 mm.

3. Effects of nonuniform field

As noted in Sec. IV, the average drift field in LUX is 180 V/cm. Here, we treat the position dependence in more detail for some relevant corrections for which using the average field is not optimal.

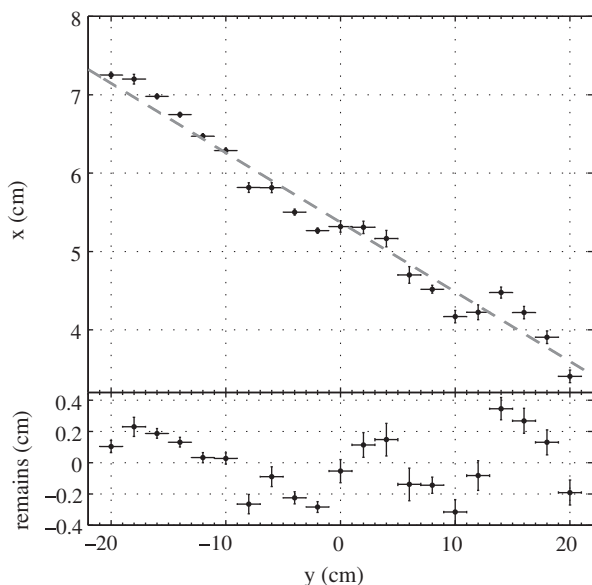


FIG. 30. *Top*: The crosses show the average reconstructed x position as a function of reconstructed y from D-D calibration data. In both cases, x and y are corrected for the effects of the nonuniform electric field. The dashed line shows a linear fit to the D-D beam direction. *Bottom*: Deviation from the linear fit in the x direction.

Due to field leakage through the cathode, the electric field in the active xenon is not perfectly uniform, resulting in a small radial component. Figure 8 (left) in Ref. [23] shows the reconstructed positions of events from ^{83}mKr calibrations. The nonaxial component to the field results in the electron cloud from an event being pushed inward, away from the wall of the detector. The inward radial pull is on average $0.15 \text{ mm}/\mu\text{s}$ (toward the center of the chamber). The position reconstruction is performed using the S2 signal, resulting in events being assigned an xy position based on where in the liquid the electrons were extracted. Consequently, the calculated xy position is closer to the center than the true xy position, with a vertical dependence arising from the field's depth dependence.

To understand this phenomenon, the detector electric field was modeled in two dimensions using COMSOL version 4.3b. The model was assumed to be asymmetric and included the grids and insulators with their proper voltages and dimensions. The predicted field strength is shown in Fig. 8 (right) in Ref. [23]. At high radius, the field lines become nonparallel, leading to events being reconstructed at smaller radius than the true event position. Along the vertical axis of the detector ($x = y = 0$), the field varies from 120 V/cm just above the cathode to 220 V/cm just below the gate, leading to an average of 180 V/cm. This variation has a negligible effect on the light and/or charge yields from low-energy nuclear or electron recoils. Furthermore, any such variation in ER response over the fiducial volume is averaged out in the measurement of the tritium ER band.

The distribution of ^{83}mKr is distributed uniformly throughout the detector, allowing it to be used to produce a mapping of reconstructed xy positions to real xy positions. The detector is cut into $30 \mu\text{s}$ slices in drift time, and each slice is then further segmented into 60 sections in the polar directions. The first 30 min of any injection are ignored to ensure enough time for uniform mixing. The remaining events in each section are placed into 600 uniform radial bins, and the average radius in each bin is used to calculate the radial correction. This scheme enforces the reconstructed positions to be radially uniform within each drift time slice. Figure 31 shows the effects of this radial correction for data with drift times between 200 and $300 \mu\text{s}$. Effects of nonuniform fields on drift time are found to be negligible.

The field nonuniformity also has a small effect on the pulse-area corrections. Because position-dependent pulse-area corrections are applied based on ^{83}mKr data, they are affected by the nonuniformity of the field. In principle, this would not be a problem if the light and charge yields for events in the WIMP search region had the same dependence with field as ^{83}mKr events. Unfortunately, however, at lower energies, the light and charge yields are less sensitive to field. This results in a systematic uncertainty on the S2 corrections that can be calculated through comparison to tritium data and stands at 4% at the center of the detector.

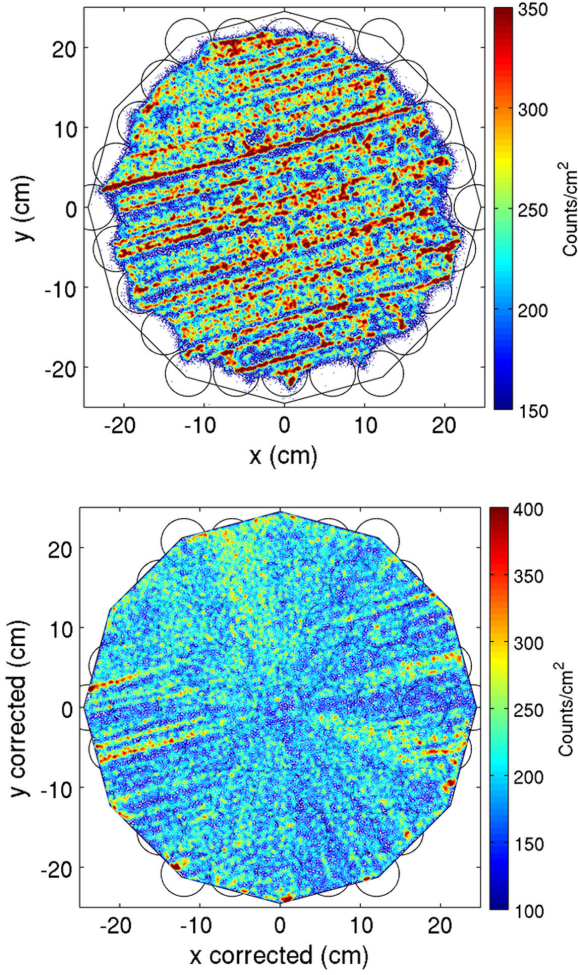


FIG. 31. *Top*: uncorrected xy positions for $^{83\text{m}}\text{Kr}$ data with drift times between 200 and 300 μs . *Bottom*: The same data, after the correction to xy position has been applied.

The effect on the S1 signal is less than 2% at any location in the detector.

D. Electron-recoil response

1. Combined energy model for electron recoils

Single-scatter events in the TPC are interpreted with the combined energy model [64]

$$E_{\text{total}} = W \cdot (n_e + n_\gamma) = W \cdot \left(\frac{S1}{g_1} + \frac{S2}{g_2} \right), \quad (5)$$

where g_1 and g_2 represent gain factors in units of phd/quantum that convert S1 and S2 signals to electron number (n_e) and photon number (n_γ), respectively. W is the energy scale factor of LXe in units of eV/quantum, g_1 is the product of the average photon collection efficiency and the average QE of the PMTs, while g_2 is the product of the electron extraction efficiency at the liquid-gas surface (ϵ_{ee}) and the single-electron size. For ER events in LUX, a constant W value of 13.7 eV/quantum is assumed [65].

TABLE II. Table of sources used in the analysis of Fig. 33.

Source	E (keV)	Type	Origin
^{127}Xe	5.3	L-shell x ray	Run 3 data
$^{83\text{m}}\text{Kr}$	41.55	IC	Internal calibration source
$^{131\text{m}}\text{Xe}$	163.9	IC	Early Run 3 data
^{127}Xe	208.3	L-shell x ray + γ	Run 3 data
^{127}Xe	236.1	K-shell x ray + γ	Run 3 data
$^{129\text{m}}\text{Xe}$	236.1	IC	early Run 3 data
^{127}Xe	409	K-shell x ray + γ	Run 3 data
^{214}Bi	609	γ	Detector background
^{137}Cs	661.6	γ	External calibration source

All source data were collected at 180 V/cm. The $^{129\text{m}}\text{Xe}$ decay and one of the ^{127}Xe processes completely overlap at 236.1 keV. IC = internal conversion.

The gain factors g_1 and g_2 may be determined by observing two or more ER line sources of known energy in which the average light and charge yields differ. g_1 and g_2 are then fixed by requiring that E_{total} , computed with Eq. (5), reproduces the true energy of each source. In ER events, the average yields vary with energy and electric field due to changes in the average recombination of ionization electrons with Xe^+ ions.

The nine sources listed in Table II were used to extract values for g_1 and g_2 in LUX. A scatter plot of $S1_c$ vs $S2_c$ for these data is shown in Fig. 32. A strong anticorrelation between S1 and S2 is apparent in each line due to recombination fluctuations. The data are fitted with a rotated two-dimensional Gaussian to determine $\langle S1 \rangle$ and $\langle S2 \rangle$ for each line source. To reduce the dependence of the result on the data selection, each fit has data selected within two Gaussian widths of the mean, as determined by the initial fit. Variation in the S2 signal and S1 signal due to

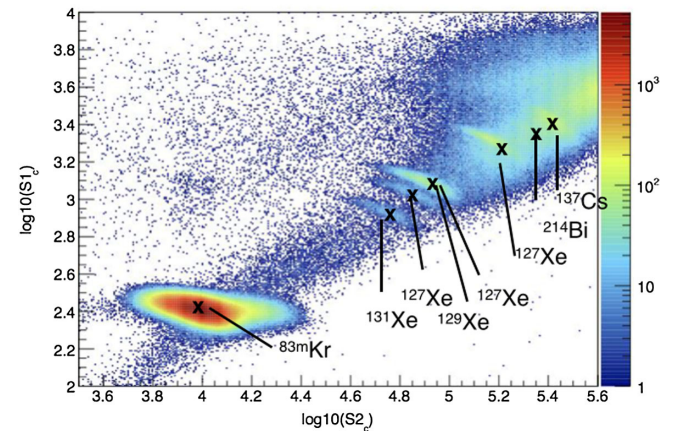


FIG. 32. $S1_c$ vs $S2_c$ for a compilation of LUX line source data. $^{83\text{m}}\text{Kr}$ and ^{137}Cs data were collected during dedicated calibration runs. All other lines were present in the low background WIMP search data. ^{127}Xe [66], ^{129}Xe , and ^{131}Xe were only present early in Run 3 due to their cosmogenic origin. The anticorrelation between S1 and S2 is due to recombination.

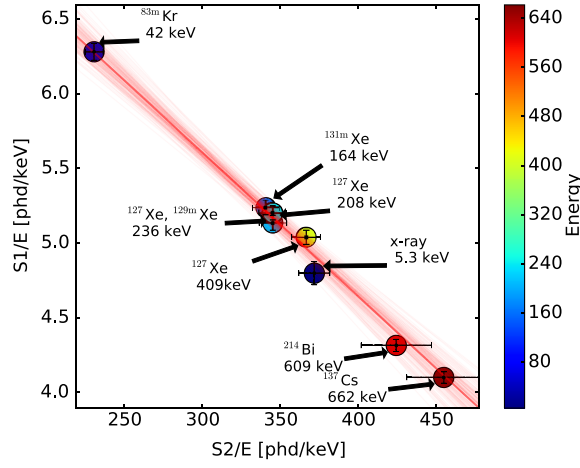


FIG. 33. Extraction of the gain factors g_1 and g_2 , adhering to the method described in the text. Each point represents a single line source listed in Table II.

PMT saturation and single-electron size variation are included as systematic errors in each fit.

The values of g_1 and g_2 were extracted by plotting $x = \langle S_2 \rangle / E$ vs $y = \langle S_1 \rangle / E$ for each source and electric field value, as shown in Fig. 33. A linear fit, $y = mx + b$, was then performed to the eight points, where $m = -g_1/g_2$ and $b = g_1/W$. Values of $g_1 = 0.117 \pm 0.003$ phd/photon and $g_2 = 12.1 \pm 0.8$ phd/electron were found, implying $\epsilon_{ee} = 0.491 \pm 0.032$. The errors were determined by toy Monte Carlo where each of the eight points was varied within its error. The extracted value for g_1 is in good agreement with the results of the optical model described in Sec. IV B.

The observed value of ϵ_{ee} may be compared to previous measurements if the electric field values above and below the LUX liquid surface are known. Uncertainty in the location of the liquid level between the gate grid and the anode leads to uncertainty in the electric field, such that a precise comparison is not possible. The COMSOL field model described in Sec. V C 3 indicates that the observed value of ϵ_{ee} is consistent with measurements from Refs. [55,56] for a liquid level 3.6 mm above the gate grid. Such a liquid level is consistent with expectations based upon the design of LUX and uncertainties in the thermal expansion coefficients of the TPC insulators. These results have been further validated by fitting tritium data to the tritium beta spectrum, finding $g_1 = 0.115 \pm 0.005$ phd/photon and $g_2 = 12.1 \pm 0.9$ phd/electron, implying $\epsilon_{ee} = 0.509 \pm 0.038$ [67].

E. Electron recoils: Implications for NEST ER model

The electron-recoil model is based upon the high-statistics tritium data sets described in Ref. [67]. The light and charge yields in simulations were tuned to reproduce the tritium data. A comparison between the tuned simulation and the tritium data is shown in Fig. 34.

As shown in Ref. [67], NEST version 0.98 for ER from 2013 [68] disagrees by $\sim 10\%$ at the lowest and highest

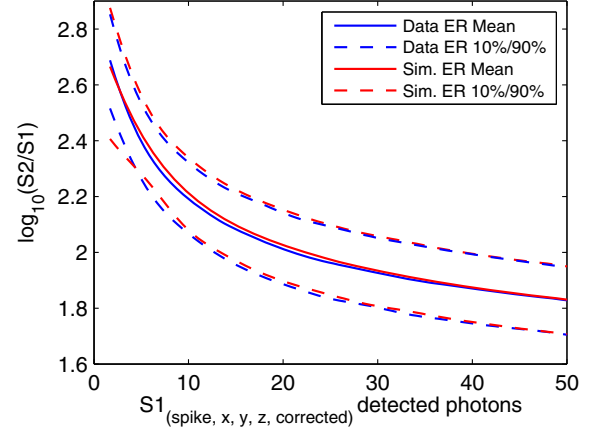


FIG. 34. The mean and width of the ER band as determined from Gaussian fits to the tritium population in S1 slices are depicted in blue, with a tuned MC based on the fundamental yields extracted using the known beta energy spectrum, at the same discrete points in S1, in red. The latter is based on a fit to the yields, but not to the band, and it is thus in a sense a prediction.

energies of interest. As thoroughly discussed in Ref. [38], NEST uses the Thomas-Imel model of recombination, but here, an extra energy dependence was added to force agreement with the high-statistics LUX tritium data for all S1 and S2 yields. This is done on a purely empirical basis, although a first-principles physics justification is to be explored in the future that will, by extension, also explore the ER S2/S1 band behavior. This was not required in previous analyses and still is not required for NR analysis, as seen next.

F. Nuclear-recoil response

1. D-D neutron calibration

Monoenergetic neutrons from a D-D fusion source were used to measure the response of the LUX detector in S_{1c} and S_{2c} to nuclear recoils in the ranges 1.1–74 and 0.7–74 keV, respectively [69].

The ratio of the ionization to scintillation signal is used to discriminate between nuclear and electron recoils in liquid xenon TPCs. Neutrons from the D-D source are used to calibrate the nuclear-recoil band over the S1 range used for the WIMP search analysis. Double-scatter nuclear-recoil events are used to kinematically reconstruct the energy of the first recoil to determine the response vs true recoil energy (see Refs. [69,70] for details). Subsequently, a simulated nuclear-recoil band is compared to the D-D calibration data to demonstrate the consistency of the nuclear-recoil signal model used to generate S1 and S2 PDFs for the WIMP search profile likelihood ratio analysis, described in Sec. VI E.

The nuclear-recoil band was measured using single-scatter events from the D-D calibration data set. An S2 threshold at 165 phd was applied on the raw S2 area before

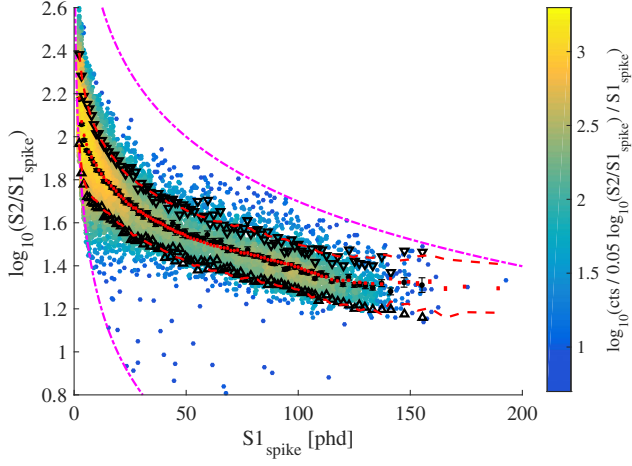


FIG. 35. The measured events from the D-D exposure defining the nuclear recoil. There are 19249 events remaining after all cuts. The black data points are the Gaussian fit mean values for each S1 spike bin. The red data points are the corresponding Gaussian fit mean values for the simulated nuclear-recoil band. The black triangles and red dashed line indicate the 90% one-sided limits from data and simulation, respectively. The lower magenta dot-dashed line indicates the S2 threshold at 165 phd, and the upper magenta dot-dashed line shows the applied upper cut for S2 ($S2 < 5000$ phd). The lower density of points in the center reflect the differential cross section behavior for 2.45 MeV neutrons incident on a xenon target. Error bars on the black and red data points are statistical only and are generally too small to see.

position correction for consistency with the WIMP search (Sec. VIB 2). The simulations were produced using LUXSIM/GEANT4 and the Lindhard-based NEST model described in Sec. VF 2 and were passed through the same LUX data-processing framework as the D-D calibration data. The same cuts and analysis used for nuclear-recoil band data were applied to the resulting reduced simulation waveforms. Figure 35 shows the nuclear-recoil band for the full range of the D-D calibration data, after the analysis cuts. The cuts applied are described in Ref. [69], but the S1 range is extended from 50 to 200 phd. This shows the data up to the recoil energy spectrum end point at 74 keV_{nr}. The simulated nuclear-recoil band is consistent with D-D calibration data within the systematic uncertainty intrinsic to the simulation process.

Figure 36 shows the same events, applying the same cuts, but now as a function of combined energy in keV_{ee}, and as a function of keV_{nr}, where keV indicates kilo-electronvolts and nr indicates the energy scale for nuclear recoils. Explicitly, the keV_{ee} energy scale is calculated through Eq. (5), where the gains during the D-D calibration time period are $g_1 = 0.115 \pm 0.004$ phd/photon and $g_2 = 11.5 \pm 0.9$ phd/electron [69]. In turn, the keV_{nr} is calculated using

$$E[\text{keV}_{\text{nr}}] = \frac{E[\text{keV}_{\text{ee}}]}{L[\text{keV}_{\text{nr}}]}, \quad (6)$$

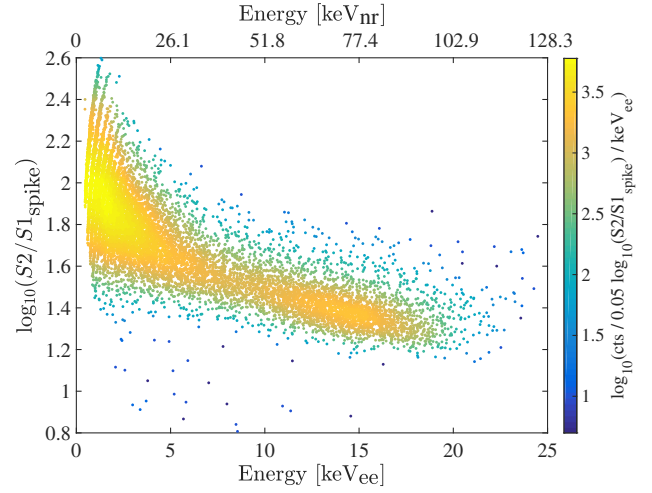


FIG. 36. The measured events defining the nuclear-recoil band are shown in the scatter plot as a function of energy in keV_{ee} and keV_{nr}. They are the same 19249 events, after analysis cuts, shown in Fig. 35. The density scale is different from the previous nuclear-recoil band plot due to normalization by bin width.

where L is Lindhard's factor [71], described in Sec. VF 2. Again, this plot shows data up to the recoil energy spectrum end point at 74 keV_{nr}. The nonzero width of the vertical bands of events at low energy is due to corrections for S1 spike overlap in the per-channel waveforms, as well as to 3D position-based detector corrections.

2. Modeling nuclear recoils in liquid xenon

A model of nuclear-recoil response for use in simulations and WIMP searches has been built and constrained following the techniques used in Ref. [39]. An energy deposit E_0 in the liquid is distributed between the formation of excitons and of electron-ion pairs. Some energy is additionally lost as an unmeasurable dissipation of heat. The process is modeled with a modified Platzman equation for rare gases [64], to which an efficiency factor, L , is applied to account for the energy lost to atomic motion rather than the detectable electronic channels,

$$E_0 = \frac{(N_{\text{ex}} + N_{\text{i}})W}{L}, \quad (7)$$

where W is the average energy required to produce a quantum (either an exciton or ion) in the liquid. Again, $W = 13.7$ eV is assumed.

Three different stages determine how the generated quanta are divided into the final light (S1) and charge (S2) signals. At the first stage, a fixed exciton-to-ion ratio $N_{\text{ex}}/N_{\text{i}}$ is defined, that determines how much of the energy initially goes to excitation as opposed to ionization. At the second stage, the ion-electron pairs recombine with probability r to produce further excitons. At the third stage, biexcitonic collisions cause some excitons to deexcite

either through heat or through Penning ionization. This effect is modeled by multiplying the resulting number of photons by a fraction f_l and allowing some fraction P of the quenched excitons to become ions. At the end of these three stages, the leftover excitons deexcite to produce scintillation photons, while the electrons that escape recombination contribute to the charge signal. The final equations for the number of photons (n_{ph}) and electrons (n_e) produced by an energy deposition E_0 are

$$n_{ph} = L \times f_l \times \frac{E_0}{W} [1 - A(1 - r)], \quad (8)$$

$$n_e = L \times \frac{E_0}{W} [B(1 - A)(1 - r) + A], \quad (9)$$

where $A = 1/(1 + N_{ex}/N_i)$ and $B = P(1 - f_l)$. The expressions for L , f_l , and r are described in detail below.

The Lindhard factor.—A factor L determines the fraction of energy in a nuclear-recoil event that goes into scintillation and ionization, rather than atomic motion. In a single Xe-Xe collision, this fraction is given by $s_e/(s_e + s_n)$, where s_e and s_n are the electronic and nuclear stopping powers (energy loss per unit distance). L represents this fraction for the entire cascade of collisions in an ionization event. The model given by Lindhard's theory [71] is adopted, with

$$L = \frac{kg(\epsilon)}{1 + kg(\epsilon)}, \quad (10)$$

where k is a proportionality constant between s_e and the velocity of the recoiling nucleus and ϵ is a dimensionless energy scale equal to $11.5(E_{nr}/\text{keV})Z^{-7/3}$. The function g models the ratio of electronic to nuclear stopping powers under the Thomas-Fermi approximation and is given by

$$g(\epsilon) = 3\epsilon^{0.15} + 0.7\epsilon^{0.6} + \epsilon. \quad (11)$$

The parameter k is treated as a free parameter for the fitting of this model to the nuclear-recoil data. (For ER, $L = 1$ fits all measurements to date.)

In addition to Lindhard's theory, an alternative model has been investigated that gives more favorable signal strength at energies below 2 keV. For this, the Thomas-Fermi approximation is replaced with the Ziegler *et al.* parametrization of nuclear stopping power, as described in Ref. [72]. Ziegler's expression is calculated using the universal screening function. The reduced stopping powers (given in terms of the dimensionless energy) in this model are given by

$$s_e(\epsilon) = 0.166\epsilon^{0.5}, \quad (12)$$

$$s_n(\epsilon_Z) = \frac{\ln(1 + 1.1383\epsilon_Z)}{2[\epsilon_Z + 0.01321\epsilon_Z^{0.21226} + 0.19593\epsilon_Z^{0.5}]}, \quad (13)$$

where $\epsilon_Z = 1.068\epsilon$. The slight difference in energy scales is due to different assumed screening lengths in the calculation of the dimensionless energy. Following Ref. [72], an additional prefactor α is introduced, which multiplies the entire expression to account for the cascade of collisions generated by a single initial nuclear recoil; thus, $\alpha > 1$. The final expression for the L -factor under this alternative model is

$$L = \alpha \frac{s_e}{s_e + s_n}. \quad (14)$$

The Lindhard model [Eq. (10)] is used in the signal model to set the WIMP-nucleon cross section limit, while the alternative [Eq. (14)] is used only to demonstrate the effect of more optimistic assumptions in yields at low energies.

Recombination.—The probability of recombination, r , is calculated using the Thomas-Imel box model [73], which gives

$$r = 1 - \frac{\ln(1 + N_i\varsigma)}{N_i\varsigma}. \quad (15)$$

The energy dependence in this equation is contained in the number of ions N_i . The quantity ς is dependent on the applied electric field. However, since LUX is operated and calibrated at a constant 180 V/cm, data provide no constraint on this property; consequently, the field dependence in this work is ignored, and ς is treated as a constant parameter.

Biexcitonic collisions.—A final quenching and ionization is applied to the light signal to account for Penning effects, in which two excitons can interact to produce one exciton and one photon [74], or one photon and one electron. Both processes remove quanta from the photon signal. Following the analysis in Ref. [72], this quenching is parametrized as the fraction derived from Birks' saturation law [75]

$$f_l = \frac{1}{1 + \eta s_e}, \quad (16)$$

where η is a strength parameter and s_e is given by Eq. (12). Here, f_l represents the proportion of excitons that remain; the fraction of quenched quanta is given by $(1 - f_l)$. This expression exhibits an increased quenching effect with increasing energy, due to higher excitation density along the track of the recoiling xenon atom. Penning ionization manifests as some fraction of the collisions resulting in the release of electrons. The parameter P is introduced to model the unknown ratio of energy lost to ionization vs heat in biexcitonic processes.

Constraining the model.—To fit the model, a global likelihood is constructed that is simultaneously constrained by

measurements of the light yield, charge yield, and nuclear-recoil band mean. The yields are constrained by an analytical model, while the nuclear-recoil band requires a full Monte Carlo (MC) simulation to generate the S1 and S2 signals from which the mean $\log_{10}(S2/S1)$ is calculated in bins of S1. The global likelihood can be separated into the product

$$\mathcal{L}_{\text{global}} = \mathcal{L}_{\text{yields}} \times \mathcal{L}_{\text{band}}. \quad (17)$$

The likelihood function for light and charge yields is constructed by assuming each point is a 2D Gaussian distribution in energy and yield, with the width given by the x and y uncertainties. A joint likelihood is calculated using Eq. (4.1) in Ref. [76],

$$\mathcal{L}_{\text{yields}}(\vec{\theta}|E_i, Y_i) = \prod_{i=0}^n \int_0^\infty A_i \exp\left(\frac{-(E_i - E)^2}{2\sigma_{E_i}^2}\right) \times \exp\left(\frac{-(Y_i - \mu_Y)^2}{2\sigma_{Y_i}^2}\right) dE, \quad (18)$$

where A_i is a normalization constant, E_i is the energy of the point i , Y_i is either charge or light yield (L_y or Q_y) of point i , and μ_Y is the model prediction at the energy E . The vector $\vec{\theta}$ is the vector of free parameters in the model.

The likelihood for the band mean is constrained by the mean values of $\log_{10}(S2/S1)$ in the nuclear-recoil data, binned by S1. A full MC simulation of the NR band is performed using the model defined by $\vec{\theta}$ calculating mean values in the same S1 bins to compare to the data. The likelihood equation is given by

$$\mathcal{L}_{\text{band}}(\vec{\theta}|\overline{\log_{10}(S2/S1)}_j) = \prod_{j=0}^m B_j \exp\left(\frac{-(\overline{\log_{10}(S2/S1)}_j - \mu_{\text{sim}})^2}{2(\sigma_j^2 + \sigma_{\text{sim}}^2)}\right). \quad (19)$$

In the above, B_j is a normalization constant, $\overline{\log_{10}(S2/S1)}_j$ is the measured band mean in the j th S1 bin, and σ_j is the measured uncertainty. The values for μ_{sim} and σ_{sim} are the band mean and uncertainty calculated from the simulation. Due to the need to run the simulation repeatedly while constraining the model, the simulation is run with statistics comparable to the measured values.

To optimize this model, a Metropolis-Hastings Markov-chain Monte Carlo algorithm (MCMC) was used to produce 10,000 samples of $\mathcal{L}_{\text{global}}(\vec{\theta})$ [77]. The advantage of this method is that it produces a map of the likelihood distribution across all parameters, naturally incorporating correlations. To extract an optimal model and uncertainties, the MCMC sample was used to calculate the means and standard deviations of each of the five parameters. The results are shown in Table III. The MCMC was also used to

TABLE III. Means and standard deviations of the free parameters using both the Lindhard and the alternative model, calculated from the MCMC sample set.

Parameter	Lindhard model	Alternative model
k	0.1735 ± 0.0060	N/A
α	N/A	2.212 ± 0.081
ς	0.0168 ± 0.0020	0.0145 ± 0.0018
N_{ex}/N_i	0.482 ± 0.069	0.487 ± 0.058
η	13.2 ± 2.3	6.7 ± 3.4
P	0.111 ± 0.060	0.18 ± 0.10

construct 68% and 95% confidence bands on the model. The calculation of Q_y and L_y from the best-fit model with confidence intervals is shown in Fig. 1 of Ref. [19], and the best-fit simulation of the nuclear-recoil band is shown in Fig. 13 of Ref. [69].

G. ER and NR discrimination

The tritium data used are shown in Figs. 37(a) and 37(c). In panel (c), the 248 events below the NR mean are represented by a different marker for clarity. The ER bandwidth is represented by the dashed lines and is mostly flat for S1 signals larger than 5 phd with a small broadening below this value. The difference in $\log_{10}(S2/S1)$ between the NR and ER means, shown in solid lines in the different panels of Fig. 37, has a significant variation below 20 phd, resulting in a large dependency of the discrimination on S1.

To estimate the discrimination as a function of S1, the leaking events were sliced in 3.3 phd wide bins with the first slice centered at 3.35 phd and the last at 49.55 phd. This ensured enough events in each bin to estimate the leakage fraction—defined as the ratio between the number of leaking events and the total number of tritium events observed in each bin—while maintaining consistency with the slicing used in both the NR and ER calibrations. The results are shown by black squares in Fig. 38. One can see that the leakage fraction decreases in the first few bins but then increases with S1 up to ~ 20 phd, remaining constant thereafter. This behavior could not be observed in past detectors, which saw only the smooth increase, because it requires a lower threshold (which is driven by a high g_1). The green triangles in the same figure represent the leakage fraction obtained from a pure Gaussian extrapolation; in this case, Gaussian fits to the ER band slices are used to obtain the bandwidth as a function of the S1, which is then used to estimate the number of events that are below the NR band centroid. For the first bins, this Gaussian estimate underpredicts observed leakage due to contributions from upward S1 fluctuations of events with mean energy below the threshold; these events contain relatively smaller S2s and will thus end up well below the ER band. The greatest consistency with Gaussianity occurs at high S1. The non-Gaussian or “anomalous” leakage may be explained as

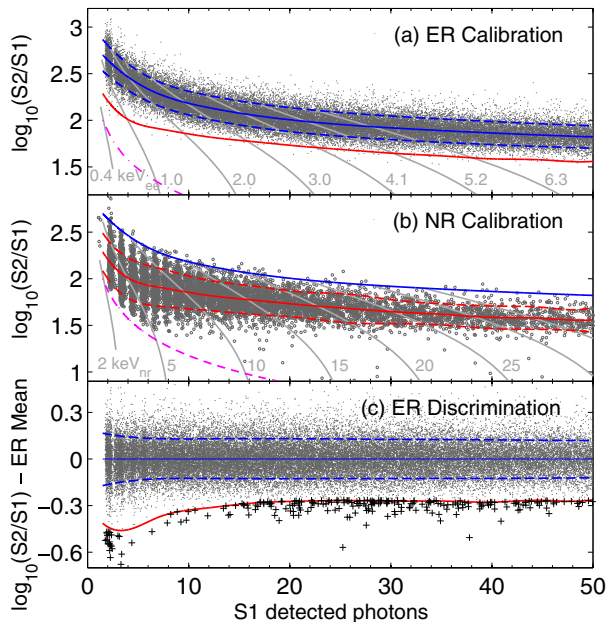


FIG. 37. Calibrations of the detector response in the fiducial volume and leakage of ER events below the NR mean. Panel (a) shows the ER (tritium) calibration, while panel (b) shows the NR calibration (obtained with monoenergetic neutrons from a D-D generator). Solid lines show band means, while dashed lines indicate the $\pm 1.28\sigma$ contours (blue for ER and red for NR). Also shown is the S2 threshold applied in the analysis (dot-dashed magenta line). Panel (c) shows $\log_{10}(S2/S1)$ normalized to the mean of the ER band as a function of the S1 for the tritium data. The 248 events that fall below the NR mean are plotted using a different marker (vertical cross) for clarity.

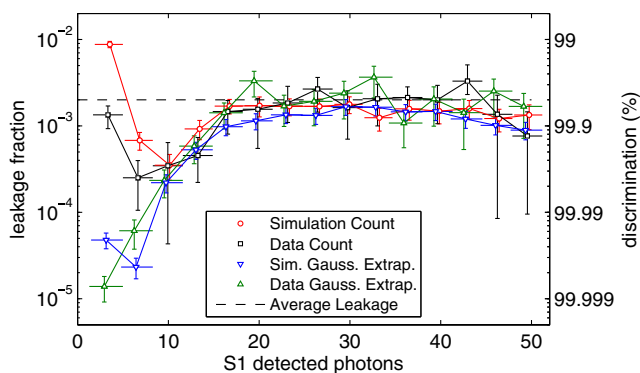


FIG. 38. Leakage (left) and discrimination (1-leakage, right) vs S1. Values obtained from a Gaussian extrapolation (blue and green triangles) are shown alongside those obtained from counting tritium events. The dashed line corresponds to the average discrimination of 99.80% in the data count. To achieve good agreement on observed leakage (data count), the g_1 and g_2 used in simulations had to be adjusted by 5%, well within uncertainty. The Monte Carlo neither underestimates nor overestimates the leakage fraction systematically, alternating between the two within error, so, for both total leakage and Gaussian only, MC (first principles) and data agree fairly well.

above and was expected based on simulation with LUXSIM and NEST in GEANT4 (red circles).

The dashed line in Fig. 38 is the weighted average of the leakage, in which the weights are the number of events in the WIMP search data for each S1 slice. The statistical uncertainty takes into account the uncertainty from the NR calibration and the binomial uncertainty from the number of the leaking events, while the systematic uncertainty is dominated by the field variations along the drift time—resulting in changes in the discrimination between different z positions in the chamber—with a smaller contribution coming from the variation of the single-electron size between the NR and ER calibrations.

Considering the number of ER events in the WIMP search run after quality and fiducial cuts are applied, the expected leakage below the NR mean is $0.7 \pm 0.1(\text{stat}) \pm 0.3(\text{syst})$ events.

VI. DATA ANALYSIS

A. Event selection: Analysis cuts

Data acquired between April 21 and September 1 2013 have been analyzed, a total of 132 calendar days, of which 101 days correspond to the data sample for the WIMP search. The remaining data samples correspond to ER and NR calibrations, xenon circulation outages, and two periods in early August following tritiated methane injections (due to a higher electron-recoil background following tritium injections, a conservative decision was taken to exclude these data from the final analysis). The final live time for the actual dark-matter search corresponds to 95 days, which accounts for exclusions due to periods of detector instability (2.5%), the trigger hold-off (1.8%), and other minor contributions (lost/corrupted files and DAQ dead time, of about 1.6%).

Single-scatter events (dubbed golden events), potentially indicative of WIMP elastic scattering off nuclei, were selected for further analysis (see Sec. III D 2). Monitoring of the detector stability was performed using slow-control (SC) parameters that could influence its response: liquid level, high voltage (HV) grid voltages and currents, outer vacuum vessel pressure, circulation flow rate, and detector pressure. All these parameters were scanned for the entire duration of the run looking for out-of-bounds periods, with excursions occurring less than 5 min apart being merged into a single unified period. Zero-length outages (a single data point outside bounds) were extended to the past and the future by the average of the update period for the corresponding SC sensor in the 10 min interval containing the excursion. Zero-length outages were not considered for the liquid level and the outer vacuum pressure, as they would likely be caused by sensor fluctuations; all other outages resulted in data being excluded from the analysis. An outage was considered to be over once the sensor returned to within nominal

bounds, with the exception of circulation failures; in these cases, the outage period was extended until the next $^{83\text{m}}\text{Kr}$ calibration, which provided updated information on the electron lifetime. Additionally, the trigger rate of each data set was visually inspected, and periods of 30 s around the rate excursions were excluded from the final analysis.

A radial fiducial cut was placed at 20 cm, driven by the leakage of decay products from ^{214}Pb (a ^{222}Rn daughter) implanted on the detector walls as discussed in Sec. VIC 2. Additionally, the height of the fiducial volume was defined to extend from 38 to 305 μs in drift time (corresponding to a total liquid height of 40.3 ± 0.2 cm) to reduce backgrounds from the PMT arrays and electrodes. Finally, the S1 and S2 thresholds (1 and 165 phd respectively, described in detail in Sec. VIB) were applied to the remaining events, as was an S1 upper limit of 50 phd.

The cleanliness of the data acquired allowed for an analysis with only a single data-quality cut, to exclude periods of a high rate of single-electron background. Small S2 pulses due to single electrons leaving the liquid surface were prevalent in the extended tails of genuine high-energy events (for up to several ms) and may have exceeded the S2 threshold either due to a fluctuation in a single-electron signal size or more commonly due to multiple electrons leaving the liquid within a few μs . These S2s may be associated with a random isolated S1 or a “fake” S1 resulting from the pulse-finder algorithm oversplitting a prior single electron and producing a false event. Excluding events with a significant area outside the (S1 and S2) signal pulses effectively removed these fake events while ensuring a large acceptance for WIMP-like events.

Figure 39 shows golden events from the WIMP search data plotted as a function of the total raw signal area

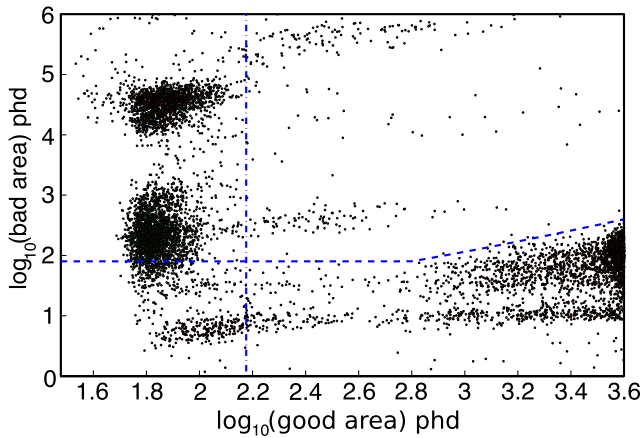


FIG. 39. A cleanliness cut applied to golden events, shown by the dashed lines, which removes periods of high-rate single-electron background, is defined as a function of extra nonsignal area contained within the 1 ms window of each event. The 165 phd S2 threshold used in the analysis is also shown (vertical dot-dashed line).

(S1 + S2 areas) and the extra (nonsignal) area contained within the same event. The dashed line indicates the threshold for the maximum allowed nonsignal area in an event, starting at 80 phd (which corresponds to ~ 3 extracted electrons) per 1 ms event window for small energy deposits and slowly increasing from this value for events with a total signal area above 630 phd. The associated acceptance was determined by slicing randomly chosen background data sets in 1 ms time windows and evaluating the total area contained within. The top plot in Fig. 40 shows the distribution of cumulative areas inside these 1 ms intervals in one of the tested data sets, with 99% of the windows containing less area than the 80 phd value used in the cut. This value was confirmed using tritium calibration data for signal areas above 400 phd (see

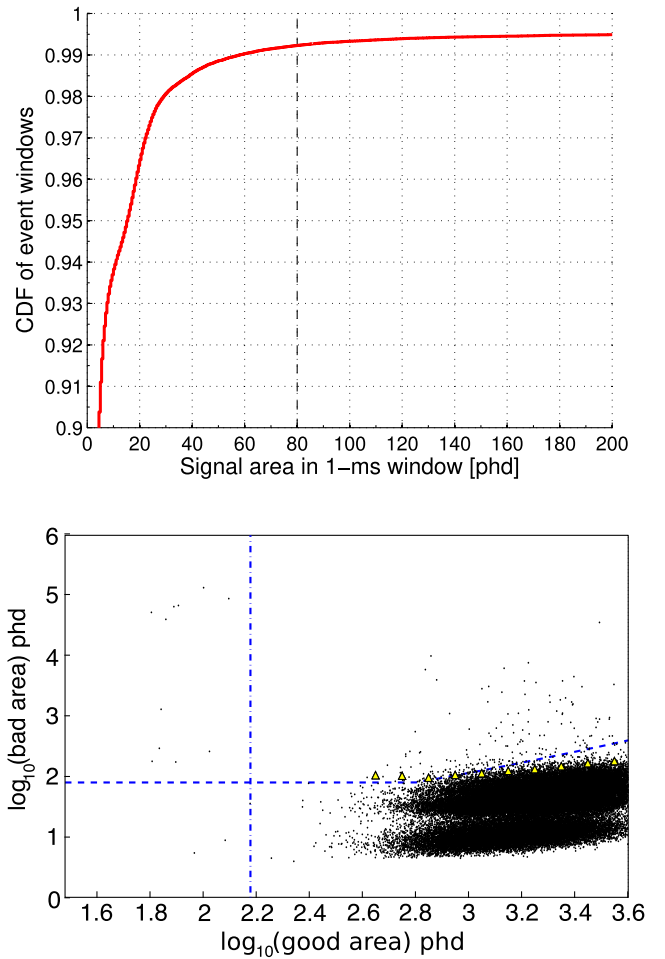


FIG. 40. Top: Cumulative number of 1 ms time window slices from a randomly picked WIMP search data set as a function of the event area contained within, showing 99% acceptance at 80 phd. Bottom: Cleanliness cuts (dashed lines) applied to a tritium calibration data set, confirming an acceptance of $\geq 99\%$ for good areas above 400 phd. Yellow triangles indicate the 99% percentile in each good area bin; the vertical dot-dashed line shows the 165 phd S2 threshold.

TABLE IV. Effect of each of the event-level cuts on the total number of acquired WIMP search triggers. Cuts have been applied sequentially.

Cut	Events remaining after cut
All triggers	93 126 362
Single scatters	5 375 716
Stability cuts	5 221 580
Fiducial volume	1 923 367
S1 range (1–50 phd)	6898
S2 threshold (> 165 phd)	840
Cleanliness cut	591

bottom plot in Fig. 40), with the cut excluding less than 1% of the events in all signal-area bins.

Table IV summarizes the effect of each of the analysis cuts and thresholds applied to the WIMP search data in the number of events acquired during Run 3.

B. Thresholds and efficiencies

1. S1 threshold

The S1 threshold is applied on the fully xyz position corrected number of spikes, the minimum of which is restricted by the two-fold coincidence requirement in the S1 definition implemented in the pulse classification module (see Sec. III D 1). As a result of position corrections, S1s smaller than two detected spikes are possible despite the two-fold coincidence, but in the original analysis [19], an additional two-phe threshold was imposed. Closer investigation of the region below this threshold did not reveal any additional noise or unexpected populations, and it was therefore decided to lower the S1 threshold to one phd for this analysis. Doing so increased the efficiency for finding a valid S1 by 15% at 3 keV_{nr} for events with a valid S2.

The fully integrated background model (described in Sec. VIC), including contributions from the cosmogenically activated ^{127}Xe that limited the upper bound of the S1 scale to 30 phe in the original analysis, allows the extension of the upper S1 limit to 50 phd. As described in Sec. VIC 2, S1s above 50 phd (outside the region of interest for the WIMP search) were used to estimate the leakage of wall events into the fiducial volume.

2. S2 threshold

The threshold on the S2 signal is applied to the uncorrected raw pulse areas and is driven by the definition of the fiducial volume and the power of the cleanliness cut to exclude events during periods of high rate single-electron background following large S2 signals.

Over a broad range of energies, events occurring close to the detector walls were found to occasionally be reconstructed to lie inside the fiducial volume. Due to incomplete

charge collection, these wall events can be mistakenly identified as WIMP-like signals. Since for small sized S2 signals these events behave in a similar manner for different S1 regions, their expected leakage into a given fiducial volume in the S1 WIMP search range of interest ($S1 < 50$ phd) may be estimated for different S2 thresholds using events from the WIMP search data with an S1 greater than 50 phd; see Sec. VIC 2 for more details on this study.

Ultimately, the S2 threshold was set at 165 phd, a compromise between the capability of the cleanliness cut to remove events in periods of high rate single-electron background (in which false S2s may be paired with coincident sphe or S1s from no-charge-collection regions of the detector) and the low efficiency for events with smaller S2s to have a valid S1.

3. Efficiencies

The efficiency to detect single scatters, which is used in the final PLR analysis, is shown by the black line in the lower panel of Fig. 1 in Ref. [19]. It was estimated using simulated NR event waveforms, generated with LUXSIM using the Lindhard parametrization to the light and charge yields obtained from the *in situ* nuclear-recoil calibrations [69]. These simulated events have been analyzed with the full data-processing framework and using the same cuts and thresholds that have been applied to the WIMP search data and show an efficiency for detecting golden events of 0.3% at the signal cutoff energy of 1.1 keV, rising to 50% at 3.3 keV and reaching 100% at around 6 keV. The drop in efficiency above 30 keV is determined by the S1 upper limit of 50 phd.

The efficiency estimated from this simulated NR data set was compared with the efficiency obtained from D-D calibration data (see Fig. 41), showing a good agreement in the relevant region of small S1 pulses. To estimate the efficiency from nuclear recoils generated by D-D neutrons, the S1 spectrum obtained from the analysis of the data, which used cuts and thresholds similar to the ones used for the WIMP search data except for a tighter fiducial cut around the beam line projection inside the detector, was compared with that obtained from a simulation of the D-D neutron beam using LUXSIM for the same number of generated neutrons. This simulation was then used to generate the energy deposits inside the xenon volume created by the nuclear recoils resulting from neutron elastic scatterings, with the S1 and S2 spectra being estimated using the same NEST version used for the D-D analysis and the PLR analysis.

To ensure that all pulse topologies occurring in the data, as well as potential losses due to the restriction to ten recorded pulses in the DPF, are covered in the efficiency estimation, an absolute DPF efficiency has been determined from 4000 NR calibration hand-scanned events, obtained with an AmBe source (see Sec. III D 2). A list of golden single-scatter events that were identified by manual

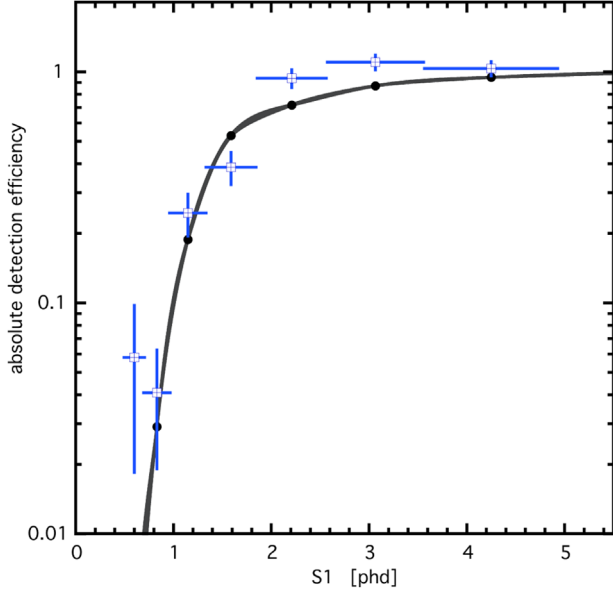


FIG. 41. Comparison of the single-scatter efficiency curve obtained from the simulation of NR events (black dots; the line is added just to guide the eye) with the efficiency obtained from analysis of D-D data (blue empty squares), in the relevant region of small S1 pulses. For larger signals ($S1 > 5$ phd), the efficiency stays flat at 100% up to the S150 phd upper limit.

scanning was compared to the output of the data-processing framework. Applying all analysis cuts, it was found that $97.5 \pm 1.7\%$ of all golden events had been successfully captured and identified by the LUX DPF. This absolute efficiency is applied as a scaling throughout the entire energy range.

C. Backgrounds

1. Random coincidence background

Apparent golden events can be generated by the random coincidence of an S1-only event and an S2-only event. The two unrelated pulses must be separated in time by less than the maximum physical drift time, with the S1 preceding the S2. S1-only events, i.e., a minimum of two PMTs recording a photoelectron within 100 ns of one another, without an accompanying S2, may be caused by energy deposition in the subcathode dead region, random coincidence of dark-count photoelectrons, or Cherenkov light (emitted, for example, in PTFE components or the window of a PMT). S2-only events are caused by low-energy tracks in an S2-live volume, i.e., the drift or extraction regions, which generate fewer than two detected S1 photons. The z position inferred from apparent drift time in these events is a uniform random variable, so they constitute a background even in large xenon TPCs where conventional single scatters from peripheral radioactivity are excluded by the fiducial volume.

To contribute to the coincidence background, an S2 must have raw area above the 165 phd analysis threshold, be reconstructed within the fiducial radius, and occur in an

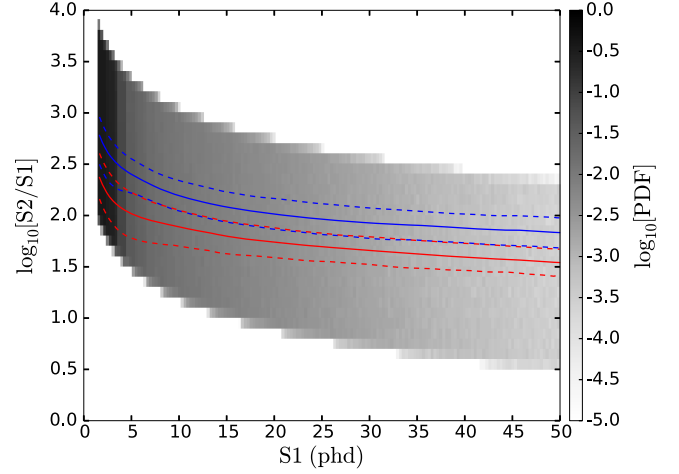


FIG. 42. Distribution in $\log_{10}(S2/S1)$ vs S1 of events caused by random coincidence of S1-only and S2-only events. The blue and red lines have the same meaning as in Fig. 37.

event window which passes the cleanliness cut and contains no preceding S1. The spectrum of S2s meeting these criteria was obtained from a representative sample of WIMP search data sets. The isolated S2 rate from threshold to 4500 phd raw area—covering the entire NR signal region—is 0.5 mHz. Beta particles and alpha-decay daughter nuclei originating on the wires of the gate electrode are a common cause of S2-only events due to the high field and the obscuration of S1 light; the characteristic short duration of gate S2s, for which the electron cloud diffuses very little in z on the way to surface, could be used to reduce this background at some penalty in efficiency, though this was not required for the LUX S1-S2 WIMP search.

The trigger efficiency for S1s rolls off to zero in the search range (below 50 phd), but, because all samples passing zero suppression are written to disk by the DAQ, the rate of isolated S1s can still be measured by applying the pulse-finder algorithm to representative WIMP search data, irrespective of the trigger. The isolated S1 rate from threshold to 50 phd raw area is 1 Hz, with a falling spectrum. Figure 42 shows the resulting distribution of background events due to coincidence S1-only and S2-only events, which leads to a prediction of 1.1 background events in the 95.0 day Run 3 reanalysis sample.

2. Nuclear-recoil and surface backgrounds

Nuclear recoils contributing to the WIMP background in LUX can be caused by elastic scattering of fast neutrons or by surface alpha decays in which the daughter nucleus recoils into the liquid xenon target. Neutron background rates are estimated in detail in Ref. [40] and contribute a negligible expectation of 0.08 events in the Run 3 reanalysis.

Wall events constitute a background to the WIMP search only when reconstructed inside the fiducial radius. Figure 43 shows the distribution of background events from 86.3 live days. The suppressed S2/S1 ratio described

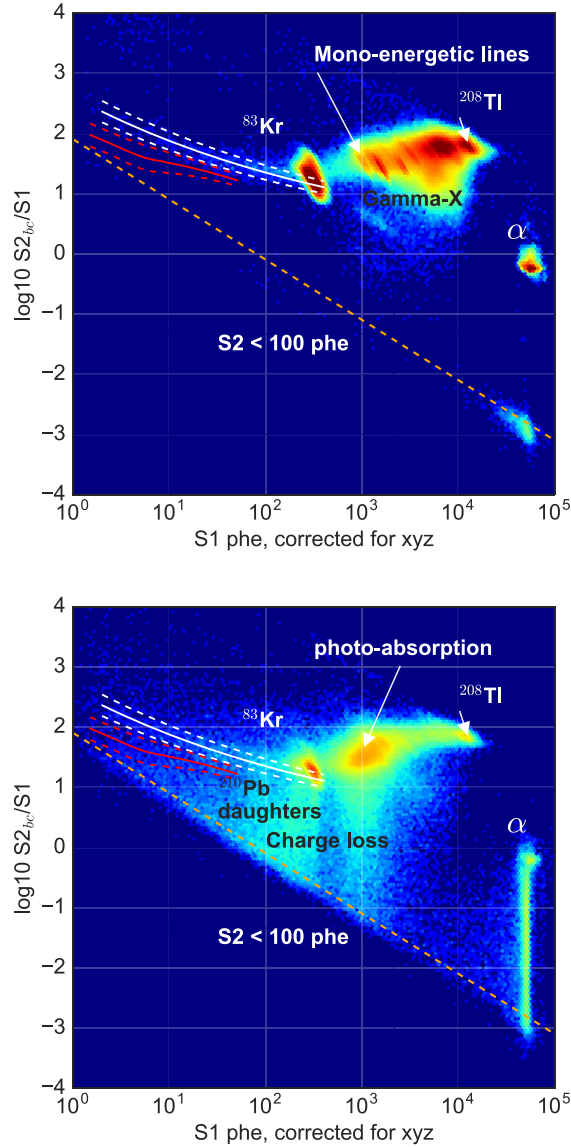


FIG. 43. Distribution in S1 vs $\log_{10}(S2_b/S1)$ of background events from 86.3 live days and a sample of ^{83m}Kr events from calibration data. Top: Events within the liquid xenon bulk defined by a radius of less than 15 cm and a drift time between 10 to 320 μs . Bottom: events near the PTFE wall, defined by the distribution of ^{210}Pb daughters. More details on the wall reconstruction can be found in Ref. [78]. The spread of events below the main peaks illustrates the incomplete charge collection for events near the wall.

above tags wall events without reference to their reconstructed radius, allowing one to model the resolution of the wall in r empirically with a high-S1, low-S2/S1 sideband away from the WIMP signal region. The low- r side of the resolved wall is well described by an S2-dependent Gaussian distribution with standard deviation

$$\sigma[\text{cm}] = \frac{1}{0.61 \log_{10}(S2_c[\text{phd}])}. \quad (20)$$

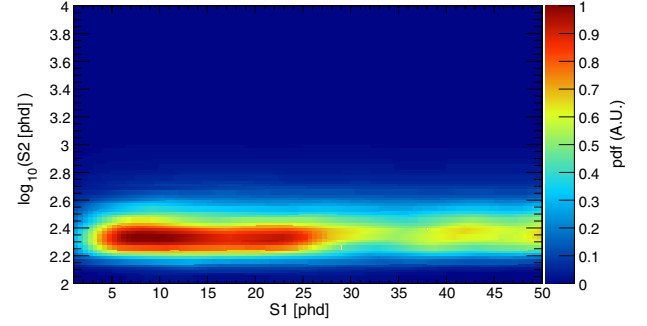


FIG. 44. Kernel-density estimated distribution, in S1 and S2, of wall background events reconstructed within the fiducial volume. The predicted count is 24 events within the 145 kg fiducial volume over 95.0 live days.

The rate and S1-S2 distribution of the wall background were derived from the data themselves, nonparametrically. X rays, Compton scatters, and ^{206}Pb daughter nuclei recoiling against undetected alpha particles all contribute to the wall population, but their signal yields are not readily simulated. Instead, an empirical wall radius was calculated, for each bin of polar angle, from the median radius of the extreme low-S2/S1 population, and kernel density estimation was applied to all events beyond this radius to construct a wall PDF in S1-S2 space. Multiplying by the Gaussian conditional distribution for r , given S2, results in the complete model for wall events in the fiducial volume, shown in Fig. 44, and a prediction of 24 ± 7 counts in the 95.0 live-day sample.

3. Electron-recoil backgrounds

Single-site electron recoils within the LUX fiducial volume are caused by β and electron-capture decays within the liquid target and by γ decays in the surrounding material. To account for systematically uncertain rates, the background model is subdivided into six populations (Table I of Ref. [19]), which are described here in descending order of the low-energy event rate.

Uranium-chain, thorium-chain, ^{40}K , and ^{60}Co nuclei in the detector construction materials emit high-energy γ rays that may, with low probability, Compton scatter once in the fiducial volume and then leave the active region. A fit of isotope activities to data in the range 0.5–3 MeV electron-equivalent energy (i.e., well above the dark-matter region of interest) gives generally good agreement [40]. Two anomalies in the fit are still under investigation, though neither has significant implications for WIMP searches; the spectrum around the 969 keV line is higher than other Th-chain activity implies, and there is an elevated continuum background in the lowest few cm of the active region. To account for the latter, the PDF for Compton scattering events was split into two source groups, with rates allowed to vary independently: one consisting of decays in the bottom PMT arrays, the surrounding support structure, and

the bottom γ -ray shield, and one consisting of decays elsewhere, principally in the top array and the sidewalls.

A spatially uniform population of beta decays, unaccompanied by other self-vetoing energy depositions in the active region, arises from two isotopes that are present in the liquid xenon target: ^{85}Kr (which is present in the atmosphere primarily due to nuclear fuel reprocessing) and ^{214}Pb (a radon daughter). A process of gas-phase chromatography was developed to reduce the former background [29]: the concentration by mass (g/g) of natural Kr in the LXe was reduced from 130×10^{-9} to 3.5×10^{-12} . Together with independent measurements of the ratio $^{85}\text{Kr}/^{\text{nat}}\text{Kr}$, this concentration implies $(0.13 \pm 0.07) 10^{-3}$ events/keV/kg/day in the region of interest with a flat energy spectrum. The rate of ^{214}Pb decays also has a spectrum that is effectively flat in the region of interest and was constrained to $0.11 \times 10^{-3} - 0.22 \times 10^{-3}$ at 90% C.L. This range is constrained from below by the rate of alpha decays from daughter ^{214}Po nuclei and from above by the total measured event rate at 300–350 keV_{ee} [40].

Electron-capture decays by ^{127}Xe and ^{37}Ar dispersed in the liquid contribute to the background of golden ER events. ^{127}Xe is a cosmogenic isotope, produced by neutron capture on ^{126}Xe . Roughly 800,000 ^{127}Xe nuclei decayed in the LUX active volume during data taking in 2013, predominantly early in the run (having a half-life of 36 days). Subsequent atomic x rays lost their energy to electrons in the bulk xenon, with the total energy determined by the shell from which the initial capture took place. Of the resulting energy lines, those at 1.1 and 5.2 keV fall in the region of interest for WIMP searches. The daughter ^{127}I nucleus emitted a γ ray at 203 or 375 keV with 43% and 57% probability, respectively. The x- and γ -ray emission were practically simultaneous, given the subnanosecond transitions and the tens-of-nanosecond scale of detector response and DAQ sampling. Only when the γ ray escaped undetected was the x-ray energy measured as a golden event, and thus a potential background. Simulation of this γ -ray inefficiency was used to predict the rate and spatial distribution of golden ^{127}Xe x-ray events, given the initial concentration of $490 \pm 95 \mu\text{Bq kg}^{-1}$ measured via the γ -ray lines.

There is some evidence for an additional line in the low-energy spectrum, consistent with a source that is uniform in position. This is thought to be due to ^{37}Ar . The fit to search data is 12 ± 8 such events in 1.4×10^4 kg days of exposure.

For all electron-recoil backgrounds, the NEST model (Sec. V E) and LUXSIM simulation of the detector and its response were used to generate MC waveforms. The same event-reconstruction software and selection cuts used for search data were applied to MC data, yielding events sampled from a simulation model that included efficiency and resolution effects. Smoothed histograms of the MC events were used to estimate the background PDFs in S1-S2 and also, for the non-spatially-uniform populations of ^{127}Xe and

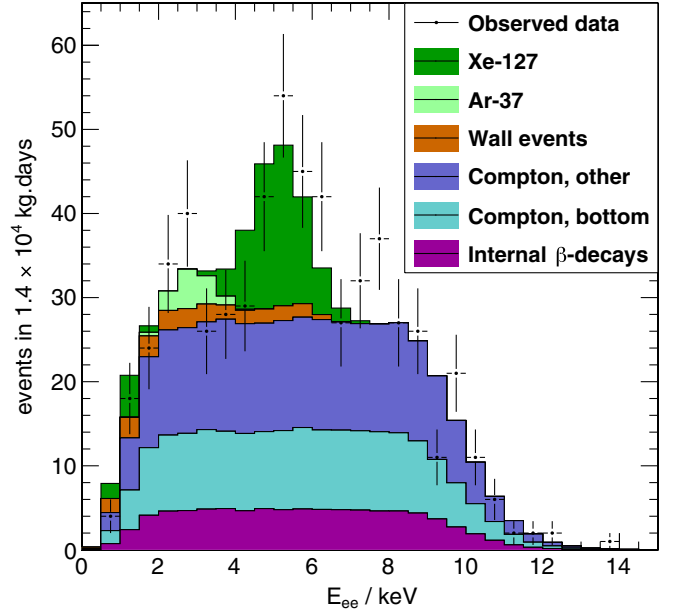


FIG. 45. Electron-equivalent-energy histogram of the LUX Run 3 search events with $S1 < 50$ phd (points) together with the best-fit model (stacked histogram), which has zero signal. The x axis is the linear, combined light-and-charge energy estimator, $E_{ee} = W \cdot (S1/g_1 + S2/g_2)$, with $W = 13.7$ eV. The roll-off in efficiency either side of 2–9 keV is primarily due to cuts on S1; search events are required to have at least two PMTs in coincidence and a maximum size of 50 phd.

Compton-scatter events, those in r - z . Figure 45 shows the background model and LUX Run 3 data as a function of recoil energy.

D. Signal model

Both spin-independent (SI) and spin-dependent (SD) WIMP-nucleon scattering models are considered in this paper. The WIMP velocity profile is derived from a standard Maxwellian velocity distribution with $v_0 = 220$ km/s and $v_{esc} = 544$ km/s as presented in Ref. [79]. The WIMP mass density, ρ_0 , is taken to be 0.3 GeV/cm^3 . For the SI case, the scattering rate is computed under the assumption that the xenon nucleons interact coherently in amplitude, thus enhancing the overall differential rate in recoil energy. In the SD case, the scalar WIMP-nucleon interaction is assumed to be suppressed, and only the axial coupling contributes to the interaction. The coherent enhancement is not present in the SD case. Typically, the axial-vector coupling constants are chosen so that the WIMP couples only to protons or only to neutrons. Xenon has an even number of protons and two naturally occurring isotopes (53% abundance) with an odd number of neutrons. Therefore, in xenon, the SD proton-only interaction is suppressed relative to neutron only. However, in the proton-only case, scattering may also occur through two-body currents (three-body interactions) where both the protons and the unpaired neutron contribute, regaining some

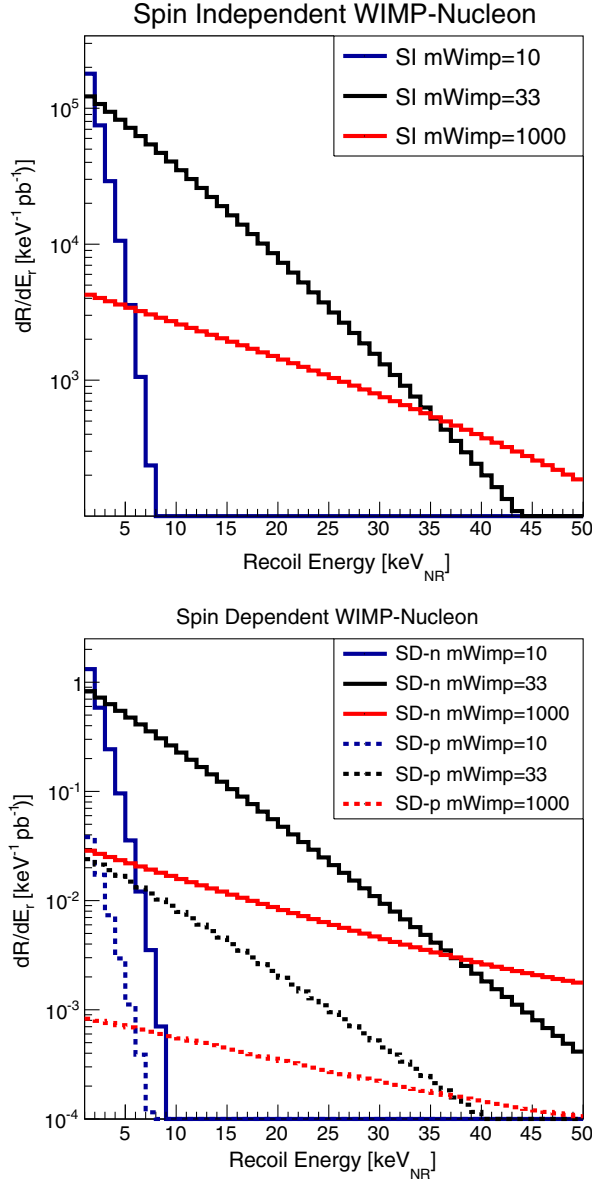


FIG. 46. Top: SI WIMP-nucleon interaction differential rate functions for WIMP mass = 10 GeV (blue), 33 GeV (black), and 1 TeV (red). Bottom: SD WIMP-neutron (solid) and WIMP-proton (dashed) interaction rate for WIMP mass = 10 GeV (blue), 33 GeV (black), and 1 TeV (red).

proton-only sensitivity. The SD model is detailed in Ref. [20] using calculations from Ref. [80]. In both the SI and the SD models, differential interaction rates are calculated as the functions of the recoil energies, as shown in Fig. 46 for some examples of WIMP masses.

Using the differential rate function for each interaction model and WIMP mass, S1 and S2 observables are generated using NEST (Sec. IV A). NEST was calibrated using *in situ* nuclear-recoil data that allowed simulation of the detector response to WIMP-nucleon interactions. Using the g_1 and g_2 obtained from the energy calibration in Sec. V D, a global fit to both the charge and light yields was

applied, using the Lindhard model [71] for the energy dependence response in nuclear-recoil interactions. These parameters are shown in Sec. V F 1 with their best-fit values and uncertainties shown in Table III. The uncertainties in these fit parameters are treated as nuisance parameter in the PLR method. For the signal models, the spatial observables, r and z , are taken to be uniformly distributed by volume. The four signal observables are combined into a four-dimensional PDF with a normalization derived from the signal efficiency (see Fig. 1 in Ref. [19]) and the total rate expected from a particular WIMP mass and interaction coupling. The signal model is defined by

$$M_{\text{sig}} = N_0 \times \epsilon(\kappa_L) \times P(\text{S1}, \log_{10}(\text{S2}), r, z), \quad (21)$$

where N_0 is the total number of events expected for a particular WIMP model, ϵ is the signal acceptance efficiency which depends on the nuisance parameter κ_L , and $P(\text{S1}, \log_{10}(\text{S2}), r, z)$. κ_L is the Lindhard factor, and $P(\text{S1}, \log_{10}(\text{S2}), r, z)$ is the four-dimensional PDF for the signal model.

To determine the effects of these nuisance parameters, the parameters were varied by $\pm 1\sigma$ from their best-fit values. Figure 47 compares the mean (top) and width (bottom) of $\log_{10}(\text{S2})$ as a function of S1 for a 50 GeV WIMP. Figure 48 shows the change of the number of events expected with our exposure for both spin-independent and -dependent WIMP models. In these plots, only the g_2 , κ_L , and Thomas-Imel box model parameter (TIB) [73] variations are shown. These distributions are different for each WIMP mass, so χ^2 , as defined by Eq. (22), is computed for each mass and parameter variation,

$$\chi^2 = \sum_{\text{S1}=1}^{\text{S1}_{\text{max}}} \frac{(Y_{\text{var}} - Y_{\text{nom}})^2 N_{\text{S1}}}{Y_{\text{nom}}}, \quad (22)$$

where Y_{nom} is the nominal value of the means or widths of the $\log_{10}(\text{S2})$ band, Y_{var} are the $\pm 1\sigma$ variation, and N_{S1} is the number of events in that particular S1 bin to account for the weights. The χ^2 values are shown in Table V for a few representative WIMP masses. The χ^2 in the means and the widths of the bands are dominated by the variation in g_2 followed by the TIB parameter. The third set of values in Table V, $\Delta(\epsilon)$, measures the change in the signal acceptance efficiency due to variation in the nuisance parameters. In the case of signal acceptance efficiency, g_2 and κ_L are the dominating nuisance parameters. Since the signal model must be regenerated during the variation of the signal model profiling, each nuisance parameter variation increases the computing time of the limits calculation by a factor of 8. We computed the exclusion limits with a few representative points with additional nuisance parameters allowed to vary. The results only fluctuate about the original by about 5%. Consequently, only κ_L and g_2 are the two nuisance parameters allowed to vary during the

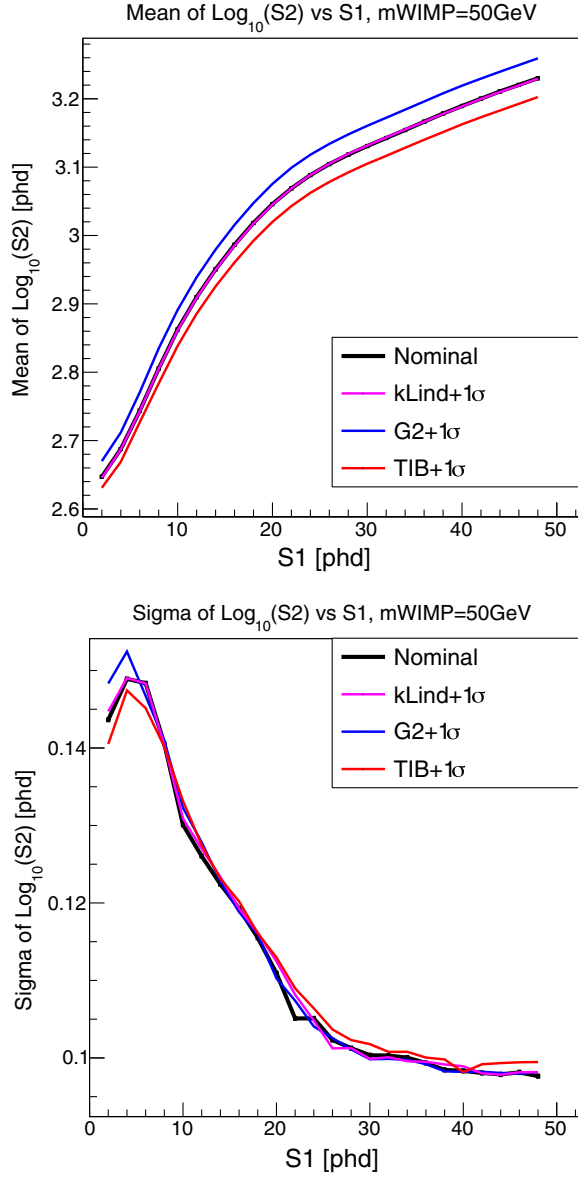


FIG. 47. Top: Mean of the $\log_{10}(S2)$ band as a function of $S1$. Bottom: Gaussian width of the $\log_{10}(S2)$ band as a function of $S1$. (Black) Nominal value of all nuisance parameters, (red) TIB varied by $+1\sigma$, (blue) g_2 varied by $+1\sigma$, and (magenta) κ_L varied by $+1\sigma$, g_1 , N_{ex}/N_i , η , and P variations are not plotted but they fall within the g_2 and TIB variations.

profile likelihood ratio analysis, where κ_L only affects the signal acceptance. For the SD analysis, the effects of structure factors were also explored. In particular, the factors $S_p(u)$ and $S_n(u)$ as defined in Sec. C of Ref. [80] for ^{129}Xe and ^{131}Xe were allowed to vary by $\pm 1\sigma$ in their theoretical uncertainties. For the WIMP-neutron interactions, $S_n(u)$, the signal strength changed by less than 7% for all masses. For the WIMP-proton

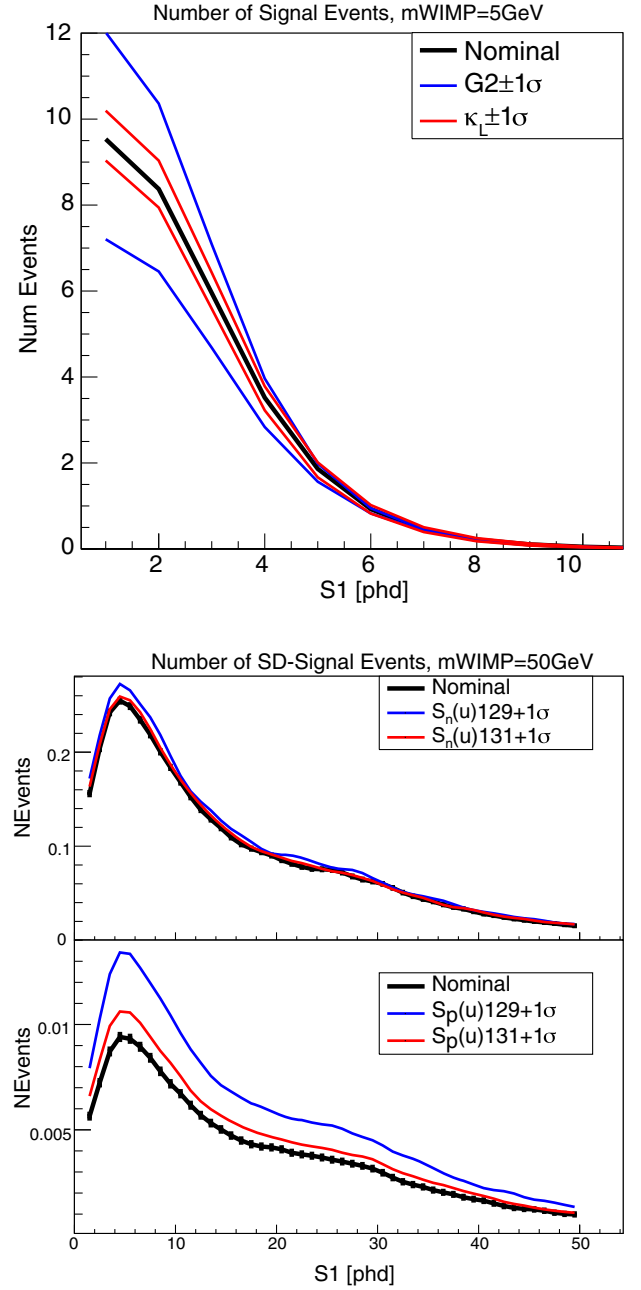


FIG. 48. Top: Number of SI WIMP-nucleon interaction events passing the analysis cuts scaled to the Run 3 exposure for a WIMP mass of 5 GeV. The black line corresponds to nominal values of all nuisance parameters, blue shows g_2 varied by $\pm 1\sigma$, and red shows κ_L varied by $\pm 1\sigma$. Bottom: The number of SD WIMP-neutron (upper) and SD WIMP-proton (lower) interaction events passing the analysis cuts scaled to the Run 3 exposure for a WIMP mass of 50 GeV. In these plots, black corresponds to nominal values of all nuisance parameters; blue shows the structure factors $S_n(u)129$ and $S_p(u)129$, as defined in Sec. C of Ref. [80], varied by $+1\sigma$; and red shows $S_n(u)131$ and $S_p(u)131$ varied by $+1\sigma$.

TABLE V. Effects of spin-independent nuisance parameter variations for various WIMP masses. Each parameter is varied by its respective $\pm 1\sigma$ uncertainty, Δ_{par} . Four representative WIMP masses and all seven possible nuisance parameter variations are tabulated here (g_1 , g_2 , and signal model parameters from Table III).

Mass	g_1	g_2	κ_L	TIB	N_{ex}/N_i	η	P
Δ_{par} (%)							
All	± 2	± 7	± 4	± 12	± 14	± 18	± 54
χ^2_{mean}							
5	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01
10	0.03	1.5	0.1	0.3	0.5	0.1	0.1
50	3.2	87	1.7	43	26	14	19
1000	0.3	9.4	0.1	6.7	1.8	2.0	2.9
χ^2_{σ}							
5	< 0.01	< 0.01	< 0.01	< 0.01	< 0.01	< 0.01	< 0.01
10	1.3	1.3	0.1	0.8	0.6	0.8	0.3
50	6.9	8.1	13	29	12	14	13
1000	1.0	2.1	1.7	2.2	2.2	2.1	1.7
$\Delta(\epsilon)$ (%)							
5	2.6	27	15	0.84	1.9	1.3	1.2
10	0.6	4.5	2.9	1.2	1.5	1.9	0.1
50	0.2	0.3	0.1	0.2	0.03	0.6	0.1
1000	0.78	0.19	1.0	0.58	0.18	2.0	0.04

interactions, $S_p(u)$, the signal strength changed by less than 40% for all masses. For this result, $S_p(u)$ and $S_n(u)$ were fixed at their central values and did not vary in the PLR.

E. Profile likelihood calculations of interaction cross section limits

A model for the LUX data and auxiliary measurements (i.e., D-D calibration and background rate estimation) is specified by a WIMP mass (m), a WIMP-nucleon cross section (e.g., σ_0 in the spin-independent analysis), and a vector of nuisance parameters (θ). The latter consists of rate parameters for each background population, together with the two signal-model parameters that have significant uncertainty: the ratio of g_2 in WIMP search vs D-D data taking and the Lindhard k parameter defined in Sec. V F. The total likelihood is a product of the likelihood of the observed data given the full model and an independent Gaussian constraint term for each element in θ , as specified in Table I of Ref. [19].

WIMP-nucleus cross section limits are computed at 31 test WIMP masses, spanning 3.5 to 20 000 GeV/ c^2 , via a PLR test. The test statistic $\lambda(\sigma)$ is the ratio of the conditional maximum likelihood at a given cross section (maximizing over the nuisance parameters with σ fixed) to the unconditional maximum likelihood. The unconditional maximum-likelihood fit to the observed data has $\sigma = 0$. Figure 49 shows the function $-\log(\lambda(\sigma))$ at an example WIMP mass. For each mass, the ROOSTATS [81] hypothesis

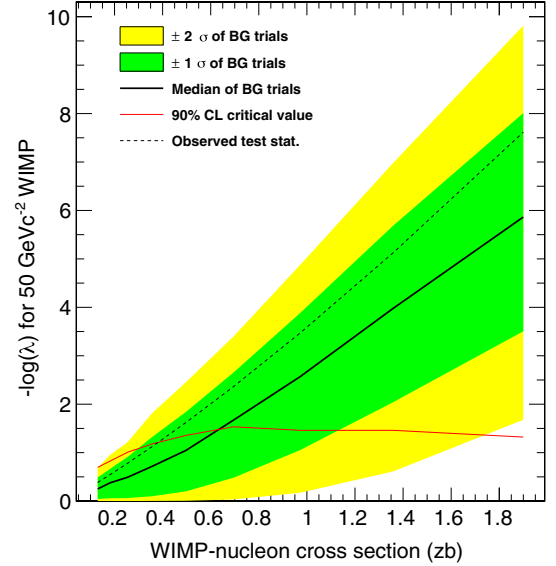


FIG. 49. Negative logarithm of the profile likelihood ratio vs spin-independent WIMP-nucleon cross section at a benchmark mass of 50 GeV/ c^2 . The critical value to exclude a cross section at 90% C.L. is in red. The raw observed test statistic is the dashed line. Given the best-fit BG-only model, the median, and the central 68% and 95% ranges are shown in black, green, and yellow, respectively.

test inverter tool is used to compute a p value for a range of cross sections and to interpolate for the observed limit, i.e., the cross section at which $(1 - p_{\text{obs}}) = \text{C.L.} = 90\%$. The same procedure is applied to MC samples from the background (BG)-only model to obtain the distribution of limits expected in the absence of signal. The raw result of the PLR test using the observed data would be a limit on the low side of its expected distribution, as seen in Fig. 49 where $-\log(\lambda(\sigma))$ lies above the median of the BG-only trials. It can be understood intuitively as the realized events fluctuating, by chance, away from the region of high signal-to-background density in the space of measured light, charge, radius, and height. This is analogous to the case of a counting experiment with background, in which the observed count happens to fall below the background-only expectation. So as not to exclude cross sections for which sensitivity is poor, a power constraint that restricts the reported limit at the median from background-only trials is applied.

The cross section limits correspond to an absolute count between 2.2–4.8 signal events in the data set, increasing monotonically with WIMP mass. In the low-mass regime, the energy spectrum of nuclear recoils is steeply falling and largely below the detection threshold in S1; the signal events that would be detected are those with upward binomial fluctuations in the number of detected S1 photons. Consequently, for a given measured S1 size, low-mass WIMP PDFs are better separated from the ER population than are high-mass ones.

VII. CONCLUSIONS AND OUTLOOK

The LUX experiment led the SI hunt for WIMP scattering until Summer 2017, and it is still currently leading the SD search for neutron coupling, thanks to progress across instrument performance, calibration sources, and data-analysis techniques. A search based on novel *in situ* calibration of PMT response and of the efficiencies and PDFs for signal and background event populations has been described. Light and charge calibrations reached down to 1 keV for both electron and nuclear recoils, based on absolute energy scales from x-ray lines and scattering kinematics, respectively. The linear, combined S1-and-S2 energy scale was shown to resolve ER lines from 5.3 to 662 keV. These in turn were used to calibrate an absolute scale of initial quanta, both for scintillation and ionization. The fiducial target mass has been estimated in two ways, one by a direct measurement of the fiducial volume and the other based on a dispersed source in the energy region of interest. After processing with an updated suite of event-reduction algorithms, 99.8% rejection of electron-recoil backgrounds was achieved based on S1 and S2 alone. Cross section limits were derived from a profile likelihood ratio test which, in addition, exploited background discrimination in spatial coordinates. The statistical derivations and motivations are detailed by Cowan *et al.* in Ref. [82]. These advances from the first LUX underground run may be applied to its subsequent, year-long run and also to future searches with liquid-xenon TPCs [83–85].

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